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February 5, 2016

### Mr. John Osolin

Remedial Project Manager Emergency and Remedial Response Division United States Environmental Protection Agency – Region 2 290 Broadway, 19<sup>th</sup> Floor New York, NY 10007-1866

### Mr. Raymond Souweha

Case Manager
New Jersey Department of Environmental Protection
401 E. State Street
Mailcode 401-05F
P.O. Box 420
Trenton, NJ 08625-0420

RE: 2015 Annual Groundwater Monitoring Report
Evor Phillips Leasing Company Superfund Site
Old Bridge Township, Middlesex County, New Jersey (**Program Interest #G000004877, EPA ID #NJD980654222**)

FILE: 19726 / 51308

Dear Mr. Osolin and Mr. Souweha:

On behalf of the Evor Phillips Leasing Company Superfund Site Settling Defendants (Group), O'Brien & Gere has prepared this 2015 Annual Groundwater Monitoring Report for the Evor Phillips Leasing Company (EPLC) Superfund Site (Site) in Old Bridge, New Jersey. The annual groundwater monitoring event was completed in March 2015, in accordance with United States Environmental Protection Agency (EPA) and New Jersey Department of Environmental Protection (NJDEP) comments received on January 21, 2015 regarding the Injection Event #1 Post-Injection Monitoring Report (submitted on November 21, 2014), and the Group's responses to those comments submitted to EPA/NJDEP on February 19, 2015. A Site Location Map is provided as Figure 1, and a Site Plan with well locations is provided as Figure 2.

The 2015 annual groundwater monitoring event was completed in conjunction with the baseline groundwater monitoring event for the second round of in-situ chemical oxidation (ISCO) injections. The locations of the ISCO treatment areas are shown on Figure 3.

The following sections of this letter present the details of the groundwater sampling event, including groundwater sampling activities and results.





### **GROUNDWATER SAMPLING ACTIVITIES**

In March 2015, a total of thirty (30) monitoring wells were sampled consistent with the 2014 annual monitoring event. A list of wells sampled and their corresponding screened intervals is included as **Table 1**. Wells were sampled for VOCs via USEPA Method 8260B.

Monitoring wells were sampled using low-flow purge methods in accordance with the approved Remedial Design Report (RDR) /Remedial Action Work Plan (RAWP) and the NJDEP Field Sampling Procedures Manual (FSPM). Purge water was containerized on site in 55-gallon steel drums. Water-level measurements were also collected from the sampled wells, and groundwater elevation data are included in **Table 2**. Groundwater elevations for the shallow aquifer are depicted in **Figure 4**.

In accordance with the approved RDR/RAWP, selected wells were also sampled for dissolved iron/chromium and total chromium/sodium via USEPA Method 6010C, total dissolved solids (TDS) via USEPA Method SM 2540C, and sulfate via USEPA Method 300.0. These samples were collected at the following wells based on proximity to the ISCO treatment areas (refer to Figure 3):

- ISCO treatment area wells (10): ISCO-MW-1 through ISCO-MW-3, ISCO-MW-5 through ISCO-MW-9, PZ-1S, IW1-BT-2
- ISCO downgradient wells (3): ISCO-MW-4, MW-14S (two sample intervals), MW-10S
- ISCO upgradient wells (3): MW-5I, IW1-DR-1, MW-11I

The analytical results for the March 2015 annual/baseline monitoring event are presented in **Table 3**, including results for quality assurance/control (QA/QC) samples (collected in accordance with the approved RDR/RAWP).

Groundwater data were validated in accordance with the approved RDR/RAWP. Data validation results are included as **Attachment 3**.

### **GROUNDWATER SAMPLING RESULTS**

### **Groundwater Flow**

Consistent with previous groundwater monitoring, shallower portions of the local aquifer are monitored via the "S" wells and the "I" wells. The deeper portions of the local aquifer are monitored via the "D" wells. Six wells (ISCO-MW-2, ISCO-MW-3, ISCO-MW-4, ISCO-MW-7, ISCO-MW-8, and ISCO-MW-9) are screened within a localized on-site perched groundwater zone (as identified by Arcadis in previous Site reports) located directly above a silty clay unit near the southern property boundary.

The shallow groundwater aquifer elevation contours are depicted in Figure 4. The two deep wells (MW-15D and MW-23D), two intermediate wells (MW-23I and WCC-1M), and the six wells screened within the localized on-site perched groundwater zone were not considered in the evaluation of the shallow groundwater elevation contours.

Consistent with historical results, shallow groundwater flow is generally toward the southwest.

### **Groundwater Quality**

2015 annual/baseline groundwater monitoring data are presented in **Table 3**. Twenty-two (22) monitoring wells screened in the shallower portion of the local aquifer (MW-6S, MW-10S, MW-14S, MW-19S, MW-23S, MW-24, MW-28, PZ-1S, IW1-BT-2, IW1-DR-1, IW-4S, WCC-1S, MW-5I, MW-9I, MW-11I, MW-23I, WCC-1M, WCC-3M, and extraction well, EW-3, and ISCO-MW-1, ISCO-MW-5, and ISCO-MW-6) and six (6) monitoring wells screened in the localized perched groundwater zone (ISCO-MW-2, ISCO-MW-3, ISCO-MW-4, ISCO-MW-7, ISCO-MW-8, and ISCO-MW-8).



MW-9) were sampled and analyzed for VOCs. Two (2) monitoring wells in the deep aquifer (MW-15D and MW-23D) were also sampled and analyzed for VOCs.

A summary of groundwater sampling results from the 2015 annual/baseline sampling event is as follows:

- Consistent with prior sampling results, VOCs with concentrations above New Jersey Groundwater Quality Standards (NJGWQS) are primarily 1,2-dichloroethane (1,2-DCA) and trichloroethene (TCE)
- Total Contaminants of Concern (COCs)¹ detected in monitoring wells at the Site and immediately downgradient from the Site were generally less than 13 ug/L
- Site COC (i.e., 1,2-DCA and TCE) concentrations were relatively unchanged in off-Site wells
- The highest total COCs were detected in three monitoring wells screened within the ISCO treatment areas, ISCO-MW-2, ISCO-MW-3, and ISCO-MW-5, which exhibited total COCs of 880 ug/L, 80 ug/L, and 200 ug/L, respectively
- The highest 1,2-DCA concentration (834 ug/L) was detected at ISCO-MW-2
- The highest TCE concentration (77 ug/L) was detected at ISCO-MW-3

The 2015 annual/baseline groundwater monitoring results are shown on Figure 5, along with estimated TCE and 1,2-DCA iso-concentration contours. Historical monitoring results are included as Attachment 1, and 2014 baseline/post-Injection Event #1 monitoring results are included as Attachment 2.

In Treatment Area 1, several wells (e.g., ISCO-MW-7, ISCO-MW-8, IW1-BT-2) exhibited total COCs 6 ug/L or less, supporting that progress toward the remedial action objectives has been made. However, groundwater results for other wells (ISCO-MW-2, ISCO-MW-3, and to a lesser degree ISCO-MW-9) continued to exhibit VOC concentrations above NJGWQS.

In Treatment Area 2, PZ-1S, ISCO-MW-1, and ISCO-MW-6 exhibited total COCs of less than 6 ug/L. As noted above, ISCO-MW-5 exhibited 200 ug/L total COCs.

Comparison of the 2015 annual/baseline results to the historical/post-round 1 results indicates that VOC concentrations in groundwater are generally either static or declining. VOC trend analyses were completed for all sampled wells and are included as **Attachment 4**.

A second round of ISCO injections to address groundwater contaminants in Treatment Areas 1 and 2 was completed June 1-11, 2015. Monitoring wells within and proximate to the treatment areas were sampled at frequencies defined in the 2015 Annual/Baseline Groundwater Monitoring Event Summary and Injection Event #2 Treatment Recommendations Summary (approved by EPA/NJDEP on May 27, 2015), to evaluate groundwater quality following the ISCO injection work. As described in that Summary, the VOCs analysis method was updated from EPA Method 8260B to Method 8260C, in accordance with the latest SW-846 methods established by EPA. This change did not materially impact the quality/usability of the data collected for this program.

In accordance with the approved RDR/RAWP, the post-injection report is being submitted to EPA/NJDEP concurrently with this 2015 Annual Groundwater Monitoring Report.

<sup>&</sup>lt;sup>1</sup> Total COCs represents the sum of those groundwater constituents above their respective New Jersey Groundwater Quality Standard (NJGWQS).



Should you have any questions regarding this submission or require additional information, please do not hesitate to contact me at (732) 638-2930.

Very truly yours, O'BRIEN & GERE ENGINEERS, INC.

**Gary Angyal, PE** Vice President

cc: EPLC Site Group

Mr. Chris Young, *de maximis, inc.* Mr. Matt Grubb, *de maximis, inc.* 

Mr. Jeffrey Levesque, O'Brien & Gere Engineers, Inc. Ms. Jessica Lehigh, O'Brien & Gere Engineers, Inc.

### **ATTACHMENTS:**

Table 1 – Monitoring Well Summary

Table 2 - March 2015 Annual/Baseline Groundwater Elevations Summary

Table 3 - March 2015 Annual/Baseline Groundwater Analytical Results

Figure 1 - Site Location Map

Figure 2 - Site Plan

Figure 3 – ISCO Monitoring Plan

Figure 4 - Groundwater Elevations - Shallow Groundwater Aquifer - March 2015

Figure 5 – 2015 Annual/Baseline Groundwater Monitoring Results

Attachment 1 – Historical Groundwater Analytical Results

Attachment 2 - 2014 Baseline/Post-Injection Event #1 Groundwater Analytical Results

Attachment 3 - Data Validation Results

Attachment 4 - Concentration Trend Graphs





# Evor Phillips Leasing Company (EPLC) Superfund Site Old Bridge, New Jersey Monitoring Well Summary Table 1

Well ID	Easting (NAD83)	Northing (NAD83)	TOC Elevation (ft MSL)	Top of Screen (ft bgs)	Bottom of Screen (ft bgs)	Top of Sample Interval (ft bgs)	Bottom of Sample Interval (ft bgs)
ISCO-MW-1	584,217.85	540,637.61	42.63	24	29	24	29
ISCO-MW-2	584,319.63	540,795.20	48.92	16	21	16	21
ISCO-MW-3	584,387.22	540,912.08	51.28	22	27	22	27
ISCO-MW-4	584,325.53	540,918.26	44.67	15	20	15	20
ISCO-MW-5	584,250.24	540,698.22	47.81	25	30	25	30
ISCO-MW-6	584,302.97	540,784.57	48.78	27	32	27	32
ISCO-MW-7	584,334.67	540,870.99	46.3	18	23	18	23
ISCO-MW-8	584,360.38	540,879.45	50.19	19	24	19	24
ISCO-MW-9	584,422.18	541,020.50	48.79	20	25	20	25
IW1-BT-2	540,925.16	584,418.94	52.39	15	35	24	29
IW1-DR-1	540,926.52	584,458.57	57.46	20	35	25	30
IW-4S	540,871.99	584,354.81	50.80	31	36	31	36
PZ-1S	540,551.93	584,158.57	44.24	20	30	22	27
MW-5I	540,691.57	584,309.75	49.74	30	40	30	35
MW-6S	540,482.53	584,118.03	43.54	17	32	22	27
MW-9I	540,610.57	584,300.26	48.40	32	42	32	37
MW-10S	540,619.21	584,165.36	45.27	15	30	23	28
MW-11I	540,543.75	584,212.88	47.92	27	37	27	32
MW-14S	540,781.83	584,184.87	32.03	3.5	18.5	S 7 D 12.5	12 17.5
MW-15D	540,495.94	584,398.81	41.88	90	100	90	95
MW-19S	540,887.95	584,582.32	56.09	19.5	35.5	28	33
MW-23S	540,625.52	583,937.55	27.89	20	30	20	25
MW-23I	540,620.38	583,935.25	27.89	50	60	55	60
MW-23D	540,630.28	583,939.84	27.95	90	100	90	95
MW-24	540,404.11	584,071.49	42.46	15	35	20	25
MW-28	541,108.40	584,474.91	49.87	15	35	23	28
WCC-1S	540,461.09	583,762.17	24.88	28	38	30	35
WCC-1M	540,452.25	583,758.98	26.42	45	55	48	53
WCC-3M	535,031.00	578,117.00	27.31	38	48	30	35
EW-3	540,428.73	584,097.72	44.38	20	65	20	25

Notes:

TOC = Top of Inner Casing
MSL = Mean Sea Level
BGS = Below Ground Surface



	TOC Elevation	Depth to Water <sup>1</sup>	GW Elevation
Well ID	(ft MSL)	(ft)	(ft MSL)
ISCO-MW-1	46.23	20.41	25.82
ISCO-MW-2	48.92	21.26	27.66
ISCO-MW-3	51.28	24.35	26.93
ISCO-MW-4	44.67	17.78	26.89
ISCO-MW-5	47.81	21.95	25.86
ISCO-MW-6	48.78	22.38	26.40
ISCO-MW-7	46.3	19.11	27.19
ISCO-MW-8	50.19	23.15	27.04
ISCO-MW-9	48.79	21.45	27.34
IW-BT-2	52.39	25.33	27.06
IW1-DR-1	57.46	30.28	27.18
IW-4S	50.80	24.28	26.52
PZ-1S	44.24	18.60	25.64
MW-5I	49.74	23.62	26.12
MW-6S	43.54	18.12	25.42
MW-9I	48.40	22.46	25.94
MW-10S	45.27	19.54	25.73
MW-11I	47.92	22.21	25.71
MW-14S	32.03	6.00	26.03
MW-15D	41.88	15.81	26.07
MW-19S	56.09	28.74	27.35
MW-23S	27.89	3.00	24.89
MW-23I	27.89	2.81	25.08
MW-23D	27.95	3.14	24.81
MW-24	42.46	17.46	25.00
MW-28	49.87	22.57	27.30
WCC-1S	24.88	1.25	23.63
WCC-1M	26.42	2.83	23.59
WCC-3M	27.31	6.02	21.29
EW-3	44.38	17.87	26.51

### Notes:

(1) Depth to water is measured in feet below top of inner casing

**GW= Groundwater** 

**TOC** = Top of Inner Casing

MSL = Mean Sea Level

**BGS = Below Ground Surface** 



Sample ID	D	FB	FB-FILTERED	FB	FB-FILTERED	ТВ	ТВ	ТВ	ISCO-MW-1	ISCO-MW-1	ISCO-MW-2	ISCO-MW-2	ISCO-MW-3	ISCO-MW-3	ISCO-MW-4	ISCO-MW-4	ISCO-MW-5
Lab Sample ID	NJ CLASS IIA GROUNDWATER QUALITY	JB89329-22	JB89329-22F	JB89329-32	JB89329-32F	JB89329-13	JB89329-23	JB89329-33	JB89329-24	JB89329-24F	JB89329-20	JB89329-20F	JB89329-14	JB89329-14F	JB89329-21	JB89329-21F	JB89329-26
Sample Date	CRITERIA (7/22/2010)	3/6/2015	3/6/2015	3/9/2015	3/9/2015	3/4/2015	3/6/2015	3/10/2015	3/9/2015	3/9/2015	3/6/2015	3/6/2015	3/6/2015	3/6/2015	3/6/2015	3/6/2015	3/9/2015
Matrix	ix ug/L	WATER	WATER	WATER	WATER	WATER	WATER	WATER	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW
Unit		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs)																	
Acetone	6000	ND (2.6)	-	ND (2.6)	-	ND (2.6)	ND (2.6)	ND (2.6)	ND (2.6)	-	98.4	-	17	-	ND (2.6)	-	ND (2.6)
Benzene	1	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	1.6	J -	ND (0.21)	-	ND (0.21)	-	0.91
Bromochloromethane	-	ND (0.49)	-	ND (0.49)	=	ND (0.49)	ND (0.49)	ND (0.49)	ND (0.49)	-	ND (2.4)	-	ND (0.49)	=	ND (0.49)	-	ND (0.49)
Bromodichloromethane	1	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (1.4)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
Bromoform Bromomethane	10	ND (0.31) ND (0.39)	-	ND (0.31) ND (0.39)	-	ND (0.31) ND (0.39)	ND (0.31) ND (0.39)	ND (0.31) ND (0.39)	ND (0.31) ND (0.39)	-	ND (1.6) 18.2	-	ND (0.31) ND (0.39)	-	ND (0.31) ND (0.39)	-	ND (0.31) ND (0.39)
2-Butanone (MEK)	300	ND (0.59)	<del>                                     </del>	ND (2.5)	-	ND (0.59)	ND (2.5)	ND (2.5)	ND (2.5)	-	65.4		ND (2.5)	_	ND (2.5)	-	ND (2.5)
Carbon disulfide	700	ND (0.50)	-	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	-	8.7	J -	ND (0.50)	-	ND (0.50)	-	ND (0.50)
Carbon tetrachloride	1	ND (0.24)	- 1	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (1.2)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)
Chlorobenzene	50	ND (0.27)	-	ND (0.27)	-	ND (0.27)	ND (0.27)	ND (0.27)	ND (0.27)	-	ND (1.4)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)
Chloroethane	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	-	ND (2.8)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)
Chloroform	70	ND (0.20)	-	ND (0.20)	-	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	-	3.2	J -	ND (0.20)	-	ND (0.20)	-	ND (0.20)
Chloromethane	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	ND (0.33)	-	3.7	J -	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Cyclohexane	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	ND (0.37)	ND (0.37)	ND (0.37)	-	ND (1.9)	-	ND (0.37)	-	ND (0.37)	-	0.45 J
1,2-Dibromo-3-chloropropane	0.02	ND (1.2)	-	ND (1.2)	-	ND (1.2)	ND (1.2)	ND (1.2)	ND (1.2)	-	ND (5.9)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)
Dibromochloromethane	1	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	ND (0.25)	-	ND (1.2)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)
1,2-Dibromoethane	0.03	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	-	ND (1.1)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)
1,2-Dichlorobenzene	600	ND (0.16)	-	ND (0.16)	-	ND (0.16)	ND (0.16)	ND (0.16)	ND (0.16)	-	ND (0.80)	-	ND (0.16)	-	ND (0.16)	-	ND (0.16)
1,3-Dichlorobenzene 1,4-Dichlorobenzene	600 75	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	-	ND (1.3) ND (1.2)	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)
Dichlorodifluoromethane	1000	ND (0.24)	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.73)	ND (0.24)	-	ND (1.2)	+ -	ND (0.24) ND (0.73)	-	ND (0.24) ND (0.73)	-	ND (0.24)
1,1-Dichloroethane	50	ND (0.75)	<del>                                     </del>	ND (0.75)	-	ND (0.75)	ND (0.75)	ND (0.35)	ND (0.75)	_	ND (1.7)	<del>                                     </del>	ND (0.75)	-	ND (0.75)	-	0.74 J
1,2-Dichloroethane	2	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	3.6	-	834	-	ND (0.30)	-	0.53 J	-	165
1,1-Dichloroethene	1	ND (0.50)	-	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	-	ND (2.5)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)
cis-1,2-Dichloroethene	70	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (1.6)	-	55	-	ND (0.33)	-	2.8
trans-1,2-Dichloroethene	100	ND (0.51)	-	ND (0.51)	-	ND (0.51)	ND (0.51)	ND (0.51)	0.84 J	-	ND (2.6)	-	1.5	-	ND (0.51)	-	0.88 J
1,2-Dichloropropane	1	ND (0.43)	-	ND (0.43)	-	ND (0.43)	ND (0.43)	ND (0.43)	ND (0.43)	-	ND (2.2)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)
cis-1,3-Dichloropropene	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (1.4)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
trans-1,3-Dichloropropene	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	ND (0.32)	ND (0.32)	ND (0.32)	-	ND (1.6)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)
1,4-Dioxane	-	ND (51)	-	ND (51)	-	ND (51)	ND (51)	ND (51)	ND (51)	-	ND (250)	-	ND (51)	-	ND (51)	-	ND (51)
Ethylbenzene	700	ND (0.40)	-	ND (0.40)	-	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	-	ND (2.0)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)
Freon 113	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	ND (0.45)	ND (0.45)	ND (0.45)	-	ND (2.2)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)
2-Hexanone	700	ND (1.7) ND (0.26)	-	ND (1.7) ND (0.26)	-	ND (1.7) ND (0.26)	ND (1.7) ND (0.26)	ND (1.7) ND (0.26)	ND (1.7)	-	ND (8.7) ND (1.3)	-	ND (1.7) ND (0.26)	-	ND (1.7) ND (0.26)	-	ND (1.7) 0.73 J
Isopropylbenzene Methyl Acetate	7000	ND (0.26) ND (3.1)	-	ND (0.26)	-	ND (0.26) ND (3.1)	ND (0.26)	ND (0.26)	ND (0.26) ND (3.1)	-	ND (1.5)	1 -	ND (0.26)	-	ND (0.26)	-	ND (3.1)
Methylrycolohexane	-	ND (0.22)	_	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	ND (0.22)	_	ND (1.1)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
Methyl Tert Butyl Ether	70	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	ND (0.26)	-	ND (1.3)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
4-Methyl-2-pentanone(MIBK)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	ND (1.1)	ND (1.1)	ND (1.1)	-	ND (5.3)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)
Methylene chloride	3	ND (0.81)	-	ND (0.81)	-	ND (0.81)	ND (0.81)	ND (0.81)	ND (0.81)	-	ND (4.1)	-	ND (0.81)	-	1.7 J	-	ND (0.81)
Styrene 1,1,2,2-Tetrachloroethane	100	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-	ND (0.26)	ND (0.26) ND (0.39)	ND (0.26)	ND (0.26)	-	ND (1.3)	-	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-	ND (0.26)
Tetrachloroethene	1	ND (0.35)	-	ND (0.35)	-	ND (0.39) ND (0.35)	ND (0.35)	ND (0.39) ND (0.35)	ND (0.39) ND (0.35)	-	25.1 ND (1.8)	-	2.3	-	ND (0.35)	-	ND (0.39) ND (0.35)
Toluene	600	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	ND (0.22)	-	ND (1.1)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,2,3-Trichlorobenzene	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	ND (0.26)	-	ND (1.3)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
1,2,4-Trichlorobenzene	9	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	ND (0.22)	-	ND (1.1)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,1,1-Trichloroethane	30	ND (0.32)	-	ND (0.32)	-	ND (0.32)	ND (0.32)	ND (0.32)	ND (0.32)	-	ND (1.6)	1 -	ND (0.32)	-	ND (0.32)	-	0.86 J
1,1,2-Trichloroethane Trichloroethene	1	ND (0.28) ND (0.25)	<del>                                     </del>	ND (0.28) ND (0.25)	-	ND (0.28) ND (0.25)	ND (0.28) ND (0.25)	ND (0.28) ND (0.25)	ND (0.28) 0.34 J	-	ND (1.4) 5.9	<del>                                     </del>	ND (0.28)	-	ND (0.28) <b>0.56</b> J	-	ND (0.28) 33.3
Trichlorofluoromethane	2000	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (1.4)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
Vinyl chloride	1	ND (0.17)	- 1	ND (0.17)	-	ND (0.17)	ND (0.17)	ND (0.17)	ND (0.17)	-	ND (0.87)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)
m,p-Xylene	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	ND (0.45)	ND (0.45)	0.55 J	-	2.5	J -	ND (0.45)	-	ND (0.45)	-	0.55 J
o-Xylene	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	ND (0.20)	ND (0.20)	0.28 J			J -	ND (0.20)	-	ND (0.20)	-	0.37 J
Xylene (total) Total VOCs	1000	ND (0.20) 0	<del>                                     </del>	ND (0.20) 0	-	ND (0.20) 0	ND (0.20) 0	ND (0.20) 0	0.83 J 5.61	-	1068.2	-	ND (0.20) 152.9	<del>                                     </del>	ND (0.20) 2.79	-	0.91 J 206.58
10141 4000	-	I 0	1 1 1	U	1	U	U	1 1	3.01	I l	1000.2	1	132.8	I	2.13	1	200.00
GC/MS Volatile TIC																	
Total TIC, Volatile	-	0	- 1	0	- 1	0	0	0	0	- 1	0	- 1	0	- 1	0	- 1	11.6 J
Total Alkanes	-	0	-	0	-	0	0	0	0	-	0	-	0	-	0	-	0
Metals Analysis																	
Chromium	70	<10	<10	<10	<10	-		-	11.8	<10	188	48.5	354	185	321		<10
Iron Sodium	300 50000	<10000	<100	<10000	<100	-	-	-	11,000	<100	1,070,000	5,320	176,000	431	13,200	306	17,900
	53000	1.3000	<u> </u>	1.0000		1	1	1 1	,000	<u> </u>	.,0.0,000	1	,000	1	.0,200	1	,500
General Chemistry																	
Solids, Total Dissolved	500000	<10000	-	<10000	-	-	-	-			2,480,000	<u> </u>	678,000	-	48,800	-	<10000
Sulfate	250000	<10000	<u>-</u>	<10000	-	-	-		118,000	-	1,980,000	-	259,000	-	41,200	-	79,200
	·								·		· ·			·	·	·	

Notes:

ND, < Not Detected Above Detection Limits

- Not Sampled

Bolded value indicates a detect above detection limits

Red bolded value indicates a detection that exceeds regulatory criteria

Sample ID	NJ CLASS IIA	ISCO-MW-5	ISCO-MW-6	ISCO-MW-6	ISCO-MW-7	ISCO-MW-7	ISCO-MW-8	ISCO-MW-8	ISCO-MW-9	ISCO-MW-9	IW1-BT-2	IW1-BT-2	IW1-DR-1	IW1-DR-1	IW-4S	IW-4S	PZ-1S
Lab Sample ID	GROUNDWATER QUALITY	JB89329-26F	JB89329-18	JB89329-18F	JB89329-19	JB89329-19F	JB89329-16	JB89329-16F	JB89329-30	JB89329-30F	JB89329-15	JB89329-15F	JB89329-10	JB89329-10F	JB89329-17	JB89329-17F	JB89329-11
Sample Date Matrix	CRITERIA (7/22/2010)	3/9/2015 GW - FILTERED	3/6/2015 GW	3/6/2015 GW - FILTERED	3/6/2015 GW	3/6/2015 GW - FILTERED	3/6/2015 GW	3/6/2015 GW - FILTERED	3/9/2015 GW	3/9/2015 GW - FILTERED	3/6/2015 GW	3/6/2015 GW - FILTERED	3/4/2015 GW	3/4/2015 GW - FILTERED	3/6/2015 GW	3/6/2015 GW - FILTERED	3/4/2015 GW
Watrix	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs)		ug/ L	ug/L	ug/L	ug/ L	ug/L	ug/L	ug/L	ug/L	ug/ L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Acetone	6000		ND (2.6)	- 1	ND (2.6)	-	16.8	- 1	ND (2.6)	-	ND (2.6)	- 1	ND (2.6)	-	ND (2.6)	-	ND (2.6)
Benzene	1	-	ND (0.21)														
Bromochloromethane	-	-	ND (0.49)														
Bromodichloromethane	1	-	ND (0.28)														
Bromoform	4	-	ND (0.31)														
Bromomethane	10	-	ND (0.39)	-	2.8	-	ND (0.39)										
2-Butanone (MEK)	300 700	-	ND (2.5) ND (0.50)	-	ND (2.5)	-	ND (2.5) ND (0.50)	-	ND (2.5)	-	ND (2.5) ND (0.50)						
Carbon disulfide Carbon tetrachloride	1	-	ND (0.24)	-	0.92 J	-	0.51		ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.30)	-	ND (0.50) ND (0.24)	-	ND (0.24)
Chlorobenzene	50	-	ND (0.27)														
Chloroethane	-	-	ND (0.56)														
Chloroform	70	-	ND (0.20)	-	1.1	-	1.4	-	ND (0.20)	-	0.27	J -	ND (0.20)	-	ND (0.20)	-	ND (0.20)
Chloromethane	-	-	ND (0.33)	-	0.96 J	-	ND (0.33)										
Cyclohexane	-		ND (0.37)	-	ND (0.37)		ND (0.37)	-	ND (0.37)								
1,2-Dibromo-3-chloropropane	0.02	-	ND (1.2)														
Dibromochloromethane	1	-	ND (0.25)														
1,2-Dibromoethane 1,2-Dichlorobenzene	0.03 600	-	ND (0.23) ND (0.16)	+	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)								
1,3-Dichlorobenzene	600	-	ND (0.16)	-	ND (0.16) ND (0.26)	-	ND (0.16)										
1,4-Dichlorobenzene	75		ND (0.24)	-	ND (0.24)												
Dichlorodifluoromethane	1000	-	ND (0.73)														
1,1-Dichloroethane	50	-	ND (0.35)														
1,2-Dichloroethane	2	-	ND (0.30)	-	2.5	-	3.1	-	ND (0.30)	-	4.4						
1,1-Dichloroethene	1	-	ND (0.50)														
cis-1,2-Dichloroethene	70	-	0.59 J	-	ND (0.33)	-	0.44 J	-	ND (0.33)	-	ND (0.33)						
trans-1,2-Dichloroethene	100	-	0.56 J ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)
1,2-Dichloropropane cis-1,3-Dichloropropene	1	-	ND (0.43) ND (0.28)														
trans-1,3-Dichloropropene	-	-	ND (0.32)														
1,4-Dioxane	-	-	ND (51)														
Ethylbenzene	700	-	ND (0.40)														
Freon 113	-	-	ND (0.45)														
2-Hexanone	-	-	ND (1.7)														
Isopropylbenzene	700	-	ND (0.26)														
Methyl Acetate Methylcyclohexane	7000	-	ND (3.1) ND (0.22)														
Methyl Tert Butyl Ether	70	-	ND (0.26)														
4-Methyl-2-pentanone(MIBK)	-	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)		ND (1.1)	-	ND (1.1)		ND (1.1)	-	ND (1.1)	-	ND (1.1)
Methylene chloride	3	-	ND (0.81)	-	1 J												
Styrene 1,1,2,2-Tetrachloroethane	100	-	ND (0.26) ND (0.39)														
Tetrachloroethene	1	-	ND (0.35)	-	0.68 J	-	0.92	J -	0.73	J -	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)
Toluene	600		ND (0.22)	-	ND (0.22)												
1,2,3-Trichlorobenzene	-	-	ND (0.26)														
1,2,4-Trichlorobenzene 1,1,1-Trichloroethane	9	-	ND (0.22) ND (0.32)														
1,1,2-Trichloroethane	3	-	ND (0.32)	-	ND (0.28)												
Trichloroethene	1	-	1.8	-	1.1	- 1	2.9	-	12.5	-	2	-	2.7	-	ND (0.25)	-	1
Trichlorofluoromethane	2000		ND (0.28)	-	ND (0.28)												
Vinyl chloride m,p-Xylene	1 -	<del>-</del> -	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)
o-Xylene	-	- +	ND (0.43)	-	ND (0.43)	-	ND (0.20)	-	ND (0.20)	-	ND (0.43)	-	ND (0.20)	- +		J -	ND (0.20)
Xylene (total)	1000	-	ND (0.20)	-	ND (0.20)	- 1	ND (0.20)	-		J -	ND (0.20)						
Total VOCs	-		2.95		10.06		25.63		13.23		2.27		3.14		0.21		6.4
GC/MS Volatile TIC																	
Total TIC, Volatile	-		0	-	0	- 1	0	- 1	0		0		0		0		10   J
Total Alkanes	-	-	0	- +	0	-	0	-	0	- +	0	- +	0	- 1	0	-	0
Metals Analysis																	
Chromium	70	<10	<10	<10			1,420		_	<10	712	808 <sup>8</sup>	<10				14.2
Iron Sodium	300 50000	214	91,600	<100	189,000	803	418,000	<100	33,000	<100	578,000	591	20,500	16,400	117,000	416	44,500
		ı L	,	ı <u>L</u>	,	1	,	1	,-50	1	,	- L	-,		,*	1	,
General Chemistry																	
Solids, Total Dissolved Sulfate	500000 250000	-	395,000	-	824,000	-	2,420,000	-	235,000	-	1,980,000	-	225,000	-		-	420,000 102,000
Ounaic	20000		218,000	<u> </u>	326,000	-	1,240,000		<10000	-	703,000	-	75,300	-	97,000	-	102,000
Notes:																	

Notes:

ND, < Not Detected Above Detection Limits

- Not Sampled

Bolded value indicates a detect above detection limits

Red bolded value indicates a detection that exceeds regulatory criteria

Page 2 of 4 2/3/2016

L	Sample ID  Lab Sample ID  NJ CLASS IIA	PZ-1S JB89329-11F	MW-51 JB89329-9	MW-5I JB89329-9F	MW-6S JB89329-25	MW-9I JB89329-7	MW-10S JB89329-37	MW-10S JB89329-37F	MW-11I JB89329-8	MW-11I JB89329-8F	MW-14SS JB89329-35	MW-14SS JB89329-35F	MW-14SD JB89329-36	MW-14SD JB89329-36F	MW-15D JB89329-4	MW-19S JB89329-6	MW-23S JB89329-28
	Sample Date CRITERIA (7/22/2010)	3/4/2015	3/4/2015	3/4/2015	3/9/2015	3/4/2015	3/10/2015	3/10/2015	3/4/2015	3/4/2015	3/10/2015	3/10/2015	3/10/2015	3/10/2015	3/4/2015	3/4/2015	3/9/2015
	Matrix ug/L	GW - FILTERED	GW	GW - FILTERED	GW	GW	GW	GW	GW	GW	GW	GW	GW	GW	GW	GW	GW
Volatila Organic Compounds (VC	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VC Acetone	6000	- 1	ND (2.6)	1 - 1	ND (2.6)	ND (2.6)	ND (2.6)	1 -	ND (2.6)	- 1	ND (2.6)	1 - 1	ND (2.6)	- 1	ND (2.6)	ND (2.6)	ND (2.6)
Benzene	1	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)
Bromochloromethane	-	-	ND (0.49)	-	ND (0.49)	ND (0.49)	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	ND (0.49)	ND (0.49)
Bromodichloromethane	1	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)
Bromoform	4	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)
Bromomethane 2-Butanone (MEK)	10 300	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	ND (0.39) ND (2.5)	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	ND (0.39) ND (2.5)	ND (0.39) ND (2.5)
Carbon disulfide	700	-	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)
Carbon tetrachloride	1	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)
Chlorobenzene	50	-	ND (0.27)	-	ND (0.27)	ND (0.27)	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	ND (0.27)	ND (0.27)
Chloroethane	-	-	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)
Chloroform Chloromethane	70	-	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	ND (0.20) ND (0.33)	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	ND (0.20) ND (0.33)	2.1 ND (0.33)
Cyclohexane		-	ND (0.37)	-	ND (0.33)	ND (0.37)	ND (0.37)	-	ND (0.33)	-	ND (0.33)		ND (0.37)		ND (0.33)	ND (0.33)	ND (0.37)
1,2-Dibromo-3-chloropropane	0.02	-	ND (1.2)	-	ND (1.2)	ND (1.2)	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	ND (1.2)	ND (1.2)
Dibromochloromethane	1	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)
1,2-Dibromoethane	0.03	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)
1,2-Dichlorobenzene 1,3-Dichlorobenzene	600 600	-	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	ND (0.16) ND (0.26)	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	ND (0.16) ND (0.26)	ND (0.16) ND (0.26)
1,3-Dichlorobenzene	75	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)
Dichlorodifluoromethane	1000	-	ND (0.73)	-	ND (0.73)	ND (0.73)	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	ND (0.73)	ND (0.73)
1,1-Dichloroethane	50	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)
1,2-Dichloroethane	2	-	2.1	-	ND (0.30)	1.4	7.6	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	18.9
1,1-Dichloroethene cis-1,2-Dichloroethene	70	-	ND (0.50) ND (0.33)	-	ND (0.50) ND (0.33)	ND (0.50) ND (0.33)	ND (0.50) <b>0.6</b>	J -	ND (0.50) ND (0.33)	-	ND (0.50) ND (0.33)	-	ND (0.50) ND (0.33)	-	0.92 ND (0.33)	J ND (0.50) 0.7	ND (0.50) J 8.2
trans-1,2-Dichloroethene	100	-	1.6	-	ND (0.51)	ND (0.51)	1.3	-	ND (0.51)	-	ND (0.51)	-	ND (0.51)	-	ND (0.51)	ND (0.51)	0.9 J
1,2-Dichloropropane	1	-	ND (0.43)	-	ND (0.43)	ND (0.43)	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	ND (0.43)	ND (0.43)
cis-1,3-Dichloropropene	-	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)
trans-1,3-Dichloropropene	-	-	ND (0.32)	-	ND (0.32)	ND (0.32)	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	ND (0.32)	ND (0.32)
1,4-Dioxane Ethylbenzene	700	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	ND (51) ND (0.40)	ND (51) ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	ND (51) 0.98	ND (51) J ND (0.40)
Freon 113	-	-	ND (0.45)	-	ND (0.45)	ND (0.45)	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	ND (0.45)	ND (0.45)
2-Hexanone	-	-	ND (1.7)	-	ND (1.7)	ND (1.7)	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	ND (1.7)	ND (1.7)
Isopropylbenzene	700	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)
Methyl Acetate Methylcyclohexane	7000	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	ND (3.1) ND (0.22)	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	ND (3.1) <b>0.42</b>	ND (3.1) J ND (0.22)
Methyl Tert Butyl Ether	70	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)
4-Methyl-2-pentanone(MIBK)	-	-	ND (1.1)	-	ND (1.1)	ND (1.1)	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	ND (1.1)	ND (1.1)
Methylene chloride Styrene	3 100	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	ND (0.81) ND (0.26)	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	ND (0.81) ND (0.26)	18.3 ND (0.26)
1,1,2,2-Tetrachloroethane	1	-	ND (0.20)	-	ND (0.39)	ND (0.39)	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.39)	-	ND (0.20)	ND (0.20)	ND (0.39)
Tetrachloroethene	1	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	0.42 J
Toluene 1,2,3-Trichlorobenzene	600	-	ND (0.22) ND (0.26)		ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)
1,2,4-Trichlorobenzene	9	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)
1,1,1-Trichloroethane	30	-	ND (0.32)	-	1.7	2.2	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	ND (0.32)	ND (0.32)
1,1,2-Trichloroethane Trichloroethene	3	-	ND (0.28) ND (0.25)		ND (0.28) ND (0.25)	ND (0.28) ND (0.25)	ND (0.28)	-	ND (0.28) 0.52	J -	ND (0.28) ND (0.25)	-	ND (0.28) ND (0.25)	-	ND (0.28) ND (0.25)	ND (0.28) 0.57	ND (0.28) J 8.8
Trichlorofluoromethane	2000		ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)
Vinyl chloride	1	-	ND (0.17)	-	ND (0.17)	ND (0.17)	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	ND (0.17)	ND (0.17)
m,p-Xylene o-Xylene		-	1.4 0.77	J -	ND (0.45) ND (0.20)	0.53 0.2	J ND (0.45) J ND (0.20)	-	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)		ND (0.45) ND (0.20)		ND (0.45) ND (0.20)	0.63 0.41	J 0.79 J J 0.34 J
Xylene (total)	1000	-	2.2	-	ND (0.20)	0.73	J ND (0.20)	-	ND (0.20)	-	ND (0.20)	<u> </u>	ND (0.20)		ND (0.20)	1	1.1
Total VOCs	-		5.9		1.7	4.33	10.6		0.52		0		0		0.92	3.67	58.72
GC/MS Volatile TIC																	
Total TIC, Volatile	-	- 1	0	- 1	0	0	0	-	0	- 1	0	- 1	0	- 1	0	45.6	J 0
Total Alkanes	-	-	0	-	0	0	0		0	-	0	-	0	-	0	12	<b>J</b> 0
Metals Analysis																	
Chromium	70	<10	68.3	<10	- 1	- 1	<10	<10	<10	<10	<10	<10	<10	<10	- 1	- 1	-
Iron	300	309	-	145	-	-	-	<100	-	<100	-	<100	-	<100	-	-	-
Sodium	50000	-	<10000	-	-	-	37,700	-	15,300	-	109,000	-	134,000	-	-	-	-
General Chemistry																	
Solids, Total Dissolved	500000	- 1	247,000	- 1	- 1	- 1	224,000	- 1	132,000		836,000		867,000	- 1	<u> </u>	- 1	- 1
Sulfate	250000	-	76,400	-	-	-	121,000	-	56,400	-	278,000	-	280,000	-	-	-	-

Notes:

ND, < Not Detected Above Detection Limits

- Not Sampled

Bolded value indicates a detect above detection limits

Red bolded value indicates a detection that exceeds regulatory criteria

Company   Comp																				
Column	s	Sample ID	NI CLASS IIA																	
Company   Comp																				
Column	Sam																			
1			ug/L																	
1	Volatile Organic Compounds (VOCs)	Unit		ug/L		ug/L		ug/L		ug/L		ug/L		ug/L		ug/L		ug/L		ug/L
Property		1	6000	5.0	1.1	5.2	ш	ND (2.6)	-1	ND (2.6)		ND (2.6)		ND (2.6)		ND (2.6)		ND (2.6)		6.6
Montandame					-		+-		$\dashv$	. ,			+	. ,	-	. ,	H	. ,	-	
Marcheller		+			+	, ,	+	. , ,	$\dashv$				+		1		H		1	` '
December   1					T		+		#						Ħ		H			
December   Company   Com	Bromoform	1	4	ND (0.31)		ND (0.31)	Ħ	ND (0.31)	T	ND (0.31)		ND (0.31)	T	ND (0.31)	T	ND (0.31)	Ħ	ND (0.31)	T	ND (0.31)
Stock and Stock	Bromomethane		10	ND (0.39)		ND (0.39)	Ħ	ND (0.39)	T	ND (0.39)		ND (0.39)		ND (0.39)		ND (0.39)		ND (0.39)		ND (0.39)
Control of Control o	2-Butanone (MEK)		300	ND (2.5)		ND (2.5)	Ħ	ND (2.5)	T	ND (2.5)		ND (2.5)		ND (2.5)		ND (2.5)		ND (2.5)		ND (2.5)
Commission	Carbon disulfide		700	ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)
Company   Comp	Carbon tetrachloride			ND (0.24)		ND (0.24)		ND (0.24)		ND (0.24)		ND (0.24)		ND (0.24)		ND (0.24)		ND (0.24)		ND (0.24)
December   70			50	. ,		. ,		. ,								, ,		. ,		
Commentment			-				$\perp$		4								Ш			
Continuemen						. ,	$\perp$	. ,	4	` '		. ,	_		J	. ,	Ш	. , ,		. , ,
1.500mm/s-shorourser   100			-		+	, ,	+	. ,	4				4		-		$\sqcup$		1	
2000   2000	,		- 0.03		+	. ,	+	` '	+	` '	_	. ,	+	, ,	$\dashv$	. ,	$\vdash$	. , ,	$\vdash$	` ′
2-500-condenders				. ,	+	. ,	+		+	. ,	_		+	. ,	$\dashv$	. ,	$\vdash$		$\vdash$	` '
12 Desirenterance			•		+		+		+				+		$\dashv$		$\vdash$		+	
13-Delinote-tourne	,			. ,	+	, ,	+	. ,	+				+		+	. ,	H	. ,	+	, ,
1.6-Districtoriesment   75	,	+			H	, ,	+	. ,	$\dashv$	, ,	7	. ,	+	, ,	$\dashv$	. ,	H	, ,	+	` '
Debte-containment					H		$\dagger \dagger$		$\dashv$				+		$\forall$		Ħ		H	
12-001-001-001-001-001-001-001-001-001-0	Dichlorodifluoromethane	1		ND (0.73)		ND (0.73)	Ħ	ND (0.73)	T	ND (0.73)			T		T	ND (0.73)	Ħ	ND (0.73)	T	ND (0.73)
1.0-Octoberhene	1,1-Dichloroethane		50	ND (0.35)		ND (0.35)		ND (0.35)		ND (0.35)		ND (0.35)		ND (0.35)		ND (0.35)		ND (0.35)		ND (0.35)
Earl-2 Detrovemente   78	1,2-Dichloroethane		2	1.2		ND (0.30)		ND (0.30)		ND (0.30)		ND (0.30)		33.3		ND (0.30)		ND (0.30)		0.52
Table   100   ND (0.51)   ND (0.52)   ND	1,1-Dichloroethene			ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)		ND (0.50)			J		J	ND (0.50)
1. DO-Dishopropropropropropropropropropropropropro																				
2013-13-Orderlosprogree   -       NO (0.28)   NO (0.				. ,		, ,			J				J	, ,		. ,		. ,		. ,
Trans-1-5 Chickropopene					$\perp$		$\perp$		4				_		_		Ш			
1-4 Decision					+		+	` '	4				_		$\perp$		Н		-	
Flytherane					+	, ,	+	. ,	4	, ,			_	, ,	$\perp$	, ,	Н	. ,	-	. ,
From 113					+		+		+				+		-		H		-	
2-Hearone		-			+		+		+		-		+		H		H		$\vdash$	
Supproprietation   Topic   ND (0.26)   ND (0.27)   ND (0.28)   N					+	, ,	+		$^{+}$	` '	1		+	, ,	Ħ	, ,	H	, ,	+	` '
Methylacetate   7000					+		+		$^{+}$		1		+		Ħ		H		+	
Methyl Tarbuyl Effer					T	. ,	+		#						Ħ		H			
#### Methylsee pelicone   Mol.   Mol.	Methylcyclohexane		-	ND (0.22)	T	ND (0.22)	+	ND (0.22)	#	ND (0.22)		ND (0.22)		ND (0.22)	Ħ	ND (0.22)	H	ND (0.22)		ND (0.22)
Methylene chionide   3			70				J	. , ,		. ,		. ,		, ,				. ,		
Syren   100			-				Ш					. ,		, ,			Ш			
1,1,2,2-Fishchroroethane					+		+		4			, ,	_		$\perp$		Н		-	, ,
Introchrochene		-			+		+		+		-		+		H		H		$\vdash$	ND (0.26)
12.2-Trichlorobenzene			1		T		T		T		J		$\top$		T		H		H	ND (0.35)
12.4-Inchlorobenzene   9	Toluene		600	ND (0.22)		ND (0.22)		ND (0.22)	T	ND (0.22)		ND (0.22)		ND (0.22)		ND (0.22)	Ħ	ND (0.22)	T	ND (0.22)
11.1-inchloroesthane   30			-		П		П		1		I						П			
1,1,2-Trichloroethane 3 ND (0.28) ND					+		+		4				+		$\vdash \downarrow$		$\sqcup$		$\vdash$	
Trichloroethene	,,				+		+		+		+		+		+		H		Н	
Trichlorofluoromethane	• •	-	1		J		+	. ,	J	. ,	7		+	. ,	$\dashv$		J	. ,	J	. ,
m.p-Xylene	Trichlorofluoromethane		2000	,	Ħ		Ţ		╛			ND (0.28)	ᆂ				Ħ		L	ND (0.28)
C-Xylene				,	П				I					, ,		, ,	П	. ,		, ,
Xylene (total)   1000					J		+I		4				_		$\sqcup \downarrow$		Ц		$\vdash$	
Total VOCs	•				J	. ,	+	. ,	+		_		+	` '	J		$\vdash$		$\vdash$	
Common			-		$\forall$		$\forall$		$\dashv$				+		Ť		$\forall$		+	
Total TIC, Volatile								L											•	L
Total Alkanes - 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0																				
Metals Analysis Chromium 70					Ш		Ш		1			-			J		Ш		Щ	
Chromium   70   -   -   -   -   -   -   -   -   -	I otal Alkanes		-	0	Ш	0		U		U		0		U		0		0		Ü
Chromium   70   -   -   -   -   -   -   -   -   -	Motals Analysis																			
Iron 300		1	70	-		_		- 1	-	- 1		- 1		- 1						- 1
Sodium 50000		+		-	H	-	+		$\dashv$		7		+		$\dashv$		H	-	+	-
Solids, Total Dissolved 500000				-	IT	-		<u>-</u>		-		-		-		-		-	L	-
Solids, Total Dissolved 500000																				
				-	+	-	+	-	4	-	_	-	+	-	$\dashv$	-	$\vdash \vdash$	-	$\vdash$	-
		ļ	200000		Ш									<u> </u>					L	

Notes:

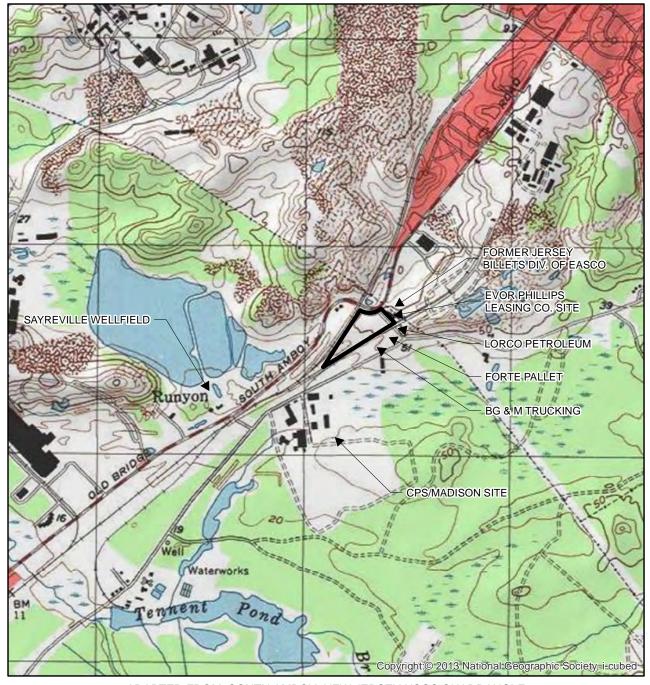
ND, < Not Detected Above Detection Limits

- Not Sampled

Bolded value indicates a detect above detection limits

Red bolded value indicates a detection that exceeds regulatory criteria





ADAPTED FROM: SOUTH AMBOY, NEW JERSEY USGS QUADRANGLE



EVOR PHILLIPS LEASING COMPANY SUPERFUND SITE OLD BRIDGE, NEW JERSEY

### SITE LOCATION



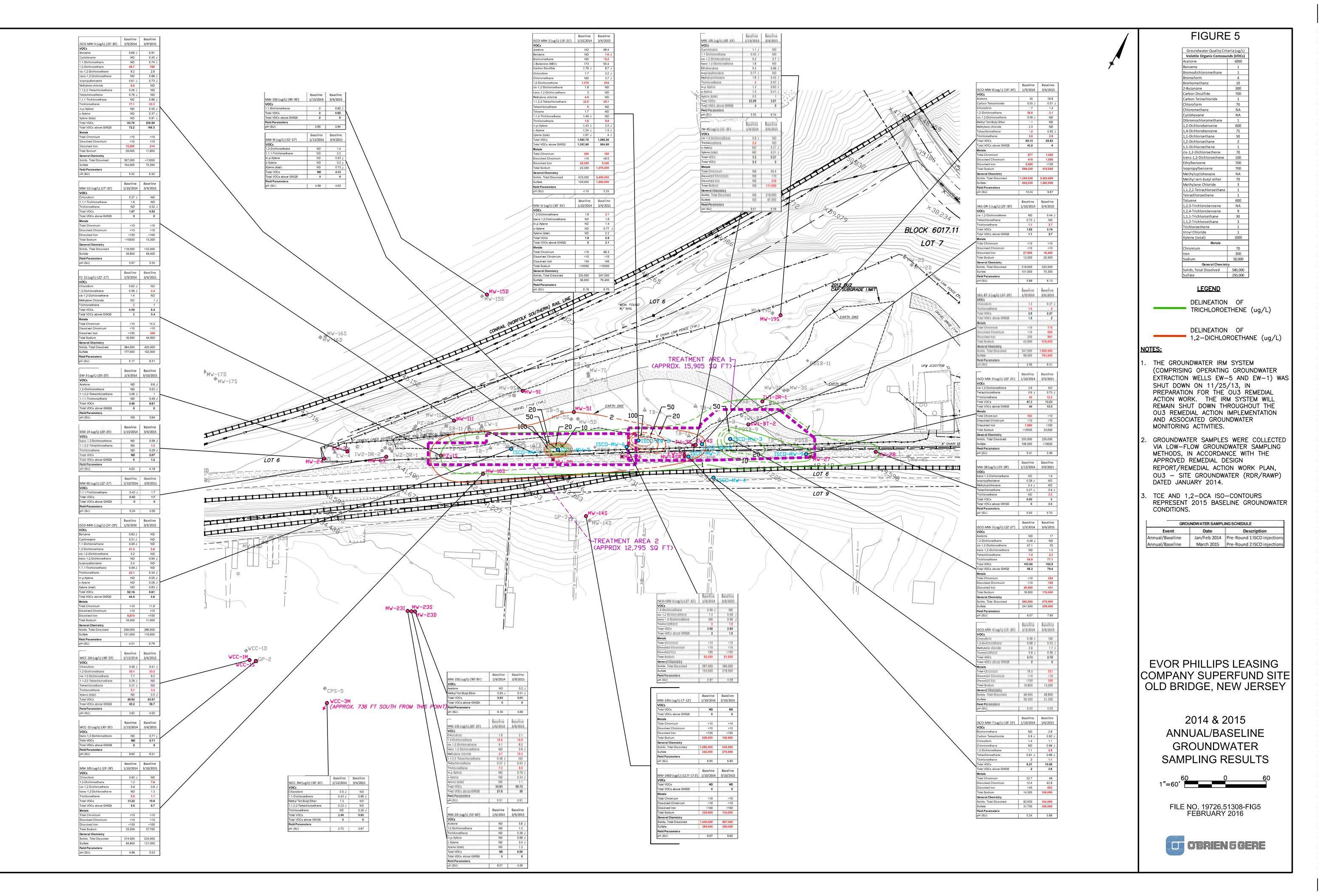


LEGEND

EXISTING CONTOUR

SANITARY SEWER MARK-OUT

FIGURE 2



# Attachment 1 Historical Groundwater Analytical Results

Sample ID	NJ CLASS IIA	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-1S	MW-4SR	MW-4SR
Sample Date		6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	6/29/2004	12/20/2004
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)			- <u></u>		<u> </u>	<i></i>	<i></i>	- J.	<i></i>		<i></i>	<u> </u>	- <u></u>		<i></i>	<u> </u>	<u> </u>	J.
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
Bromoform	4	4 U	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U [1 U]	1 U	1 U	4 U	4 U
Bromomethane	10	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U [1 U]	1 U	1 U	1.8 J	2.8
Chlorobenzene	50	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Chloroethane	-	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Chloroform	70	3.6 J	5 U	2.3 J	5 U	5 U	5 U	1.5	0.7 J	2.6 J	2.7 J	1 U	1 U	0.29 J [0.28 J]	1 U	1 U	5.3	5.9
Chloromethane	-	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Dibromochloromethane	1	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
1,2-Dichloroethane	2	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U [1 U]	1 U	1 U	54	49
1,1-Dichloroethene	1	2 U	2 U	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U [1 U]	1 U	1 U	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Ethylbenzene	700	4 U	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U [1 U]	1 U	1 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	0.8 J	1.1 J	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Methylene chloride	3	3 U	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U [1 U]	1 U	1 U	1.2 J	0.8 J
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U [20 U]	20 U	20 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	0.8 J	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	38	42
Tetrachloroethene	1	1.1	1 U	1.0	1 U	1 U	1 U	0.8	0.5 J	0.5 J	1 U	1 U	1 U	0.30 J [0.44 J]	1 U	1 U	3.5	4
Toluene	600	5 U	1.3 J	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	0.29 J	1 U [1 U]	1 U	1 U	5 U	0.6 J
1,1,1-Trichloroethane	30	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U [1 U]	1 U	1 U	3 U	3 U
Trichloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	5.4	7.2
Trichlorofluoromethane	2000	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Vinyl chloride	1	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U [1 U]	1 U	1 U	5 U	5 U
Xylene (total)	1000	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U [3 U]	3 U	3 U	5 U	5 U
Total VOCs	-	4.7 J	1.3 J	3.3 J	0.8 J	1.1 J	ND	2.3	1.2 J	3.9 J	2.7 J	ND	0.29 J	0.59 J [0.72 J]	ND	ND	109 J	112 J

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds

regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample II	NJ CLASS IIA	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-4SR	MW-5I
Sample Date	GROUNDWATER QUALITY	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/10/2012	12/20/2012	6/29/2004
Uni	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)		W5/ =		W5/ =		W5/ =		6/-	u <sub>5</sub> / =	6/-			~6/ <del>-</del>	ug/ <u>2</u>	~6/ <del>-</del>	<b>46/2</b>	-5/ -
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U [5 U]	NA
Benzene	1	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	0.6 J
Bromodichloromethane	1	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	1 U
Bromoform	4	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	4 U
Bromomethane	10	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	0.85 J [0.84 J]	NA	NA [NA]	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U [5 U]	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U [1 U]	NA
Carbon tetrachloride	1	1.1 J	2.0	1.5 J	2.0	0.6	0.9 J	2.1	3.5	2.8	1.9	0.47 J	1.0 [1.1]	1 U [0.84 J]	0.54 J	0.79 J [0.68 J]	2 U
Chlorobenzene	50	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	5 U
Chloroethane	-	5 U	5 U	5U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	5 U
Chloroform	70	4.2 J	4.6 J	4.9 J	4.8 J	2.7	3.0 J	4.8 J	5.3	5.3	2.6	1.7	3.1 [2.9]	3.0 [3.0]	2.2	3 [2.7]	5 U
Chloromethane	-	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	5 U
Dibromochloromethane	1	5 U	5 U	5 U	6 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
1,2-Dichloroethane	2	81	69	63	46	95	110	80	32	28	5.2	30	72 [67]	37 [37]	45	130 [120]	1.7 J
1,1-Dichloroethene	1	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	2 U
cis-1,2-Dichloroethene	70	0.7 J	5 U	0.5 J	5 U	1.2	0.6 J	5 U	5 U	1 U	1 U	1.1	0.36 J [0.38 J]	1 U [1 U]	0.49 J	0.35 J [0.37 J]	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA [NA]	5 U
Ethylbenzene	700	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
Methylene chloride	3	3 U	3 U	0.6 J	0.6 J	1.1	3 U	3 U	0.6 J	1 U	0.24 J	1 U	0.61 J [0.93 J]	0.37 J [0.46 J]	1 U	1.1 [0.94 J]	3 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U [20 U]	20 U [20 U]	NA	NA [NA]	100 U
1,1,2,2-Tetrachloroethane	1	38	36	45	64	20	29	42	4.8	28	2.7	9.7	10 [12]	13 [13]	9.1	20 [18]	1 U
Tetrachloroethene	1	2.4	2.9	3.2	4.5	1.7	2.5	3.1	3.1	4.9	1.6	1.0	2.1 [1.9]	0.99 J [1.0]	0.84 J	1.8 [1.5]	1 U
Toluene	600	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
1,1,1-Trichloroethane	30	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
1,1,2-Trichloroethane	3	0.6 J	3 U	0.6 J	0.6 J	0.6	0.5 J	0.4 J	3 U	1 U	1 U	0.22 J	1 U [1 U]	0.15 J [0.23 J]	0.27 J	0.32 J [1 U]	3 U
Trichloroethene	1	6.6	4.2	5.6	6.6	6.2	4.7	4.7	4.3	4.6	3.3	4.7	3.2 [3.2]	1 U [1 U]	3.4	3.1 [2.9]	0.5 J
Trichlorofluoromethane	2000	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	3.1 [3.0]	NA	NA [NA]	5 U
Vinyl chloride	1	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U [1 U]	5 U
Xylene (total)	1000	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U [3 U]	3 U [3 U]	3 U	3 U [3 U]	5 U
Total VOCs	-	134.6 J	118.7 J	124.9 J	129.1 J	129.1 J	151.2 J	137.1 J	53.6	73.6 J	17.5 J	49.9 J	92.4 J [89.4 J]	58.5 J [59.4 J]	61.84 J	160.46 J [147.09 J]	2.8 J

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-5I	MW-6S
Sample Date	GROUNDWATER QUALITY	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/10/2012	12/20/2012	6/29/2004
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U [5 U]	NA
Benzene	1	1 U	1 U	1 U	1 U	1 U	0.2	1 U	1 U	1 U	1 U	0.22 J	1 U	1 U	0.15 J	0.14 J	0.13 J [0.14 J]	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Bromoform	4	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	4 U
Bromomethane	10	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U [5 U]	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U [1 U]	NA
Carbon tetrachloride	1	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	2 U
Chlorobenzene	50	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
Chloroethane	-	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
Chloroform	70	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	2 J
Chloromethane	-	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
Dibromochloromethane	1	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
1,2-Dichloroethane	2	2.1	2 U	2 U	0.7 J	0.7 J	2.9	2 U	2 U	2 U	1.3	1.3	1.9	1.3	2.4	1.5	1.7 [1.9]	32
1,1-Dichloroethene	1	2 U	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	2 U
cis-1,2-Dichloroethene	70	5 U	5 U	5 U	5 U	5 U	0.6	5 U	5 U	5 U	1 U	1 U	1 U	1 U	0.38 J	1 U	1 U [1 U]	2.2 J
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
Ethylbenzene	700	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
Methylene chloride	3	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	27
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	NA	NA [NA]	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Toluene	600	2.7 J	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1.3	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
1,1,1-Trichloroethane	30	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
1,1,2-Trichloroethane	3	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	3 U
Trichloroethene	1	1 U	1 U	1 U	1 U	1 U	0.4	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	4.4
Trichlorofluoromethane	2000	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA [NA]	5 U
Vinyl chloride	1	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U
Xylene (total)	1000	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U [3 U]	5 U
Total VOCs	-	4.8 J	ND	ND	0.7 J	0.7 J	4.1	ND	ND	ND	1.3	2.8 J	1.9	1.3	2.9 J	1.64 J	1.83 J [2.04 J]	67.6 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-6S	MW-7I
Sample Date	GROUNDWATER QUALITY	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	7/10/2012	12/20/2012	6/29/2004
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)			-6/ -	0, -	-0/-		0/	6/-	-6/-	-6/ -	-0, -	6/ -	-0, -	6/ -			-6/ -	-87 -
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA
Benzene	1	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	0.14 J	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.66 J	1 U	1 U
Bromoform	4	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA	4 U
Bromomethane	10	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA
Carbon tetrachloride	1	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	2 U
Chlorobenzene	50	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA	5 U
Chloroethane	-	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	NA	1 U	1 U [1 U]	1 U [1 U]	NA	NA	5 U
Chloroform	70	2.1 J	2.4 J	2.0 J	1.8 J	1.8 J	1.3	0.9 J	0.9 J	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1.4	1 U	5 U
Chloromethane	-	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1.2	4.0	1 U [1 U]	1 U [1 U]	NA	NA	5 U
Dibromochloromethane	1	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
1,2-Dichloroethane	2	29	35	32	28	44	24	20	25	2 U	1 U	0.32 J	1.0	1 U [1 U]	1 U [1 U]	1 U	1 U	2 U
1,1-Dichloroethene	1	2 U	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	2 U
cis-1,2-Dichloroethene	70	2.2 J	3.5 J	8.9	13	100	18	17	12	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	0.6 J	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	NA	NA	5 U
Ethylbenzene	700	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	0.29 J	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	NA	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
Methylene chloride	3	24	18	16	12	16	10	6.2	5.8	3 U	1 U	0.26 J	0.71 J	1 U [1 U]	1 U [1 U]	1 U	1 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	NA	20 U	20 U [20 U]	2.7 J [3.3 J]	NA	NA	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	0.6 J	1 U	1.1	0.4	0.4 J	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	0.5	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.44 J	0.25 J	1 U
Toluene	600	1.8 J	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1.7	1 U	1 U [1 U]	1 U [1 U]	0.7 J	1 U	5 U
1,1,1-Trichloroethane	30	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.18 J	1 U	5 U
1,1,2-Trichloroethane	3	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	3 U
Trichloroethene	1	2.6	4.0	5.4	5.3	20	5.7	6.9	6.0	1 U	1 U	0.26 J	1 U	1 U [1 U]	1 U [1 U]	0.2 J	0.49 J	1 U
Trichlorofluoromethane	2000	5 U	5 U	5 U	1.6 J	5 U	0.5	0.8 J	5 U	5 U	1 U	NA	1 U	1 U [1 U]	1 U [1 U]	NA	NA	5 U
Vinyl chloride	1	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U
Xylene (total)	1000	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	0.74 J	3 U	3 U [3 U]	3 U [3 U]	3 U	3 U	5 U
Total VOCs	-	61.7 J	62.9 J	64.9 J	61.7 J	183.5 J	60.4	52.2 J	49.7 J	ND	ND	4.9 J	5.7 J	ND [ND]	2.7 J [3.3 J]	3.58 J	0.74 J	ND

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds

regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-7I	MW-8S	MW-8S	MW-8S
Sample Date	GROUNDWATER QUALITY	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	6/29/2004	12/20/2004	6/28/2005
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U
Bromomethane	10	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U
Chlorobenzene	50	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Chloroethane	-	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Chloroform	70	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Chloromethane	-	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Dibromochloromethane	1	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	0.9 J	5 U
1,2-Dichloroethane	2	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U
1,1-Dichloroethene	1	2 U	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Ethylbenzene	700	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Methylene chloride	3	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	100 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	1.6 J	5 U	5 U	3.1 J	5 U	0.3 U	0.4 J	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	1.8 J	0.4 J
1,1,1-Trichloroethane	30	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U
Trichloroethene	1	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	2000	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Vinyl chloride	1	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U
Xylene (total)	1000	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	5 U	1.5 J	0.9 J
Total VOCs	-	1.6 J	ND	ND	3.1 J	ND	ND	0.4 J	ND	ND	ND	ND	ND	ND	ND	ND	4.2 J	1.3 J

### Notes:

U Not Detected Above Detection Limits
-- Not Sampled
Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-8S	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I
Sample Date		12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)		<u> </u>	- U		- J,	· 0,	<u> </u>	· 0,		<u>.                                    </u>			<u></u>	<u> </u>				
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U [1 U]	0.5 J [1 U]	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U
Bromoform	4	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U [1 U]	4 U	4 U [4 U]	4 U [4 U]	4 U	4 U
Bromomethane	10	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U [1 U]	2 U	2 U [2 U]	2 U [2 U]	2 U	2 U
Chlorobenzene	50	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Chloroethane	-	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	NA	NA	NA	NA	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Chloroform	70	5 U	5 U	5 U	0.2 U	1.2 J	5 U	2.3 J	1 U	0.90 J	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Chloromethane	-	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Dibromochloromethane	1	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
1,2-Dichloroethane	2	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U [1 U]	2.1	0.9 J [0.8 J]	0.7 J [2 U]	2 U	2 U
1,1-Dichloroethene	1	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U [1 U]	2 U	2 U [2 U]	2 U [2 U]	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	5 U	5 U	0.3 U	5 U	5 U	0.3 J	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Ethylbenzene	700	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U [1 U]	4 U	4 U [4 U]	4 U [4 U]	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	0.6 J	5 U	0.3 U	5 U	5 U	5 U	1 U	NA	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Methylene chloride	3	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U [1 U]	1.5 J	0.6 J [0.6 J]	3 U [3 U]	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	NA	20 U	20 U	20 U [20 U]	100 U	100 U [100 U]	100 U [100 U]	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U
Tetrachloroethene	1	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U	1 U
Toluene	600	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	0.31 J	1 U	1 U	1 U [1 U]	5 U	2.4 J [2.3 J]	5 U [5 U]	5 U	5 U
1,1,1-Trichloroethane	30	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
1,1,2-Trichloroethane	3	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U [1 U]	3 U	3 U [3 U]	3 U [3 U]	3 U	3 U
Trichloroethene	1	0.6 J	0.4 J	0.6 J	0.4 U	0.5 J	1 U	1 U	1 U	0.43 J	1 U	0.49 J	1 U [1 U]	0.9 J	1 U [1 U]	1 U [1 U]	1 U	1 U
Trichlorofluoromethane	2000	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	NA	NA	NA	NA	0.24 J [0.24 J]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Vinyl chloride	1	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U [1 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Xylene (total)	1000	1.1 J	1.7 J	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	0.47 J	3 U [3 U]	5 U	5 U [5 U]	5 U [5 U]	5 U	5 U
Total VOCs	-	1.7 J	2.7 J	0.6 J	ND	1.7 J	ND	2.6 J	ND	1.6 J	ND	0.96 J	0.24 J [0.24 J]	4.5 J	3.9 J [3.7 J]	1.2 J [ND]	ND	ND

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds

regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-9I	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S
Sample Date	GROUNDWATER QUALITY	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/11/2012	12/20/2012	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)			-5/ <u>-</u>	~6/ <b>=</b>		w <sub>6</sub> / =	-5/ <u>-</u>	w <sub>6</sub> / =			-5/ <u>-</u>	w <sub>6</sub> / =	u <sub>5</sub> / =	~6/ <b>-</b>		W8/ =		-5/ <u>-</u>
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Benzene	1	1 U	2.5	1 U	1 U	1 U	1 U	0.22 J	1 U	1 U	1 U	0.3 J	0.31 J	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA	4 U	4 U	4 U	4 U	4 U
Bromomethane	10	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1.3 J	0.9 J	2 U	0.9 J	2 U
Chlorobenzene	50	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Chloroethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Chloroform	70	5 U	0.3	5 U	0.3 J	0.3 J	1 U	0.21 J	0.26 J	1 U	1 U	0.17 J	0.15 J	8.5	2.3 J	1.1 J	2.6 J	1.1 J
Chloromethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Dibromochloromethane	1	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,1-Dichloroethane	50	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloroethane	2	0.7 J	1.3	0.4 J	0.4 J	2 U	1 U	1 U	1.7	1 U	0.51 J	1.3	0.54 J	3.6	2.0	2 U	2 U	2 U
1,1-Dichloroethene	1	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	8.1	2.7 J	1.4 J	4.8 J	2.0 J
trans-1,2-Dichloroethene	100	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloropropane	1	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Ethylbenzene	700	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1.1 J	5 U	5 U	5 U
Methylene chloride	3	3 U	0.4 U	3 U	3 U	0.3 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	NA	NA	100 U	100 U	100 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.7 J	1 U	1 U	1 U
Tetrachloroethene	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	5 U	0.3 U	5 U	5 U	5 U	1 U	1.0	1 U	1 U	1 U	1 U	1 U	5 U	0.5 J	5 U	5 U	5 U
1,1,1-Trichloroethane	30	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.6 J	5 U	5 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U
Trichloroethene	1	1 U	1.2	1 U	1 U	1 U	1 U	1 U	1.4	1 U	1 U	0.51 J	1 U	3.3	1.0	1 U	1.1	0.8 J
Trichlorofluoromethane	2000	5 U	0.4 U	5 U	5 U	5 U	1 U	0.43 J	1 U	1 U	0.27 J	NA	NA	5 U	5 U	5 U	5 U	5 U
Vinyl chloride	1	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	0.22 J	1 U	1 U	5 U	5 U	5 U	5 U	5 U
Xylene (total)	1000	5 U	0.4 U	5 U	5 U	5 U	3 U	0.46 J	3 U	3 U	3 U	3 U	3 U	5 U	5 U	5 U	5 U	5 U
Total VOCs	-	0.7 J	5.3	0.4 J	0.7 J	0.6 J	ND	2.3 J	3.4 J	ND	1.0 J	2.28 J	1.0 J	25.4 J	11.2 J	2.5 J	9.4 J	3.9 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-10S	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I
Sample Date	GROUNDWATER QUALITY	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	7/11/2012	12/20/2012	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Benzene	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.17 J	0.088 J	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.6 J	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA	4 U	4 U	4 U	4 U	4 U
Bromomethane	10	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	0.3 U	2.8	2 U	0.7 J	1 U	1 U	1 U	0.84 J	1 U	1 U	0.29 J	2 U	2 U	2 U	2 U	2 U
Chlorobenzene	50	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Chloroethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	0.2 U	5 U	5 U	5 U	1 U	NA	1 U	NA	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Chloroform	70	5 U	1.3	6.0	5 U	4.2 J	1 U	0.51 J	0.26 J	3.9	0.29 J	1.4	1.2	5 U	5 U	5 U	5 U	5 U
Chloromethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Dibromochloromethane	1	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.31 J	1 U	5 U	5 U	5 U	5 U	5 U
1,1-Dichloroethane	50	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloroethane	2	2 U	0.3 U	1.3 J	2 U	2.3	0.9 J	16	2.9	6.2	13	1 U	1.4	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethene	1	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	2.2	3 J	5 U	2.9 J	1 U	1.4	0.77 J	11	1.9	1 U	1.3	5 U	5 U	5 U	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloropropane	1	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Ethylbenzene	700	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	0.3 U	5 U	5 U	5 U	1 U	NA	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
Methylene chloride	3	3 U	0.4 U	0.3 J	3 U	3 U	1 U	0.56 J	1 U	0.86 J	0.41 J	1 U	1 U	3 U	3 U	3 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	6.5 U	100 U	100 U	100 U	20 U	NA	20 U	20 U	20 U	NA	NA	100 U	100 U	100 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	0.4 U	1.8	1 U	1 U	1 U	1 U	1 U	0.25 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	0.4 U	0.6 J	1 U	1 U	1 U	0.22 J	1 U	0.39 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	0.8 J	5 U	5 U	5 U
1,1,1-Trichloroethane	30	5 U	0.4 U	5 U	5 U	0.4 J	1 U	0.32 J	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U
Trichloroethene	1	1 U	1.2	2.5	1 U	3.9	0.48 J	5.4	3.4	9.8	1 U	0.12 J	1.8	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	2000	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	6.5	NA	NA	5 U	5 U	5 U	5 U	1.4 J
Vinyl chloride	1	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
Xylene (total)	1000	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U	5 U	5 U	5 U	5 U	5 U
Total VOCs	-	ND	4.7	18.3 J	ND	14.4 J	1.4 J	24.4 J	7.3 J	33.2 J	22.1 J	2.6 J	6.08 J	ND	0.8 J	ND	ND	1.4 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds

regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-11I	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S
Sample Date	GROUNDWATER QUALITY	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/10/2012	12/20/2012	6/29/2004	12/21/2004	6/28/2005	12/21/2005	6/21/2006
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)								<u> </u>										
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Benzene	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1.5	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA	4 U	4 U	4 U	4 U	4 U
Bromomethane	10	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U
Chlorobenzene	50	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Chloroethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Chloroform	70	5 U	0.2 U	5 U	5 U	0.2 J	1 U	0.68 J	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
Chloromethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Dibromochloromethane	1	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,1-Dichloroethane	50	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloroethane	2	2 U	0.3 U	2 U	2 U	2 U	1 U	0.29 J	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethene	1	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloropropane	1	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Ethylbenzene	700	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.7 J
Methylene chloride	3	3 U	0.4 U	3 U	3 U	3 U	4.3	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	NA	NA	100 U	100 U	100 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	5 U	0.3 U	5 U	5 U	5 U	1 U	0.30 J	1 U	1 U	1 U	1 U	1 U	5 U	5 U	0.8 J	5 U	5 U
1,1,1-Trichloroethane	30	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U
Trichloroethene	1	1 U	0.4 U	1 U	1 U	1 U	0.42 J	2.0	1 U	1 U	1 U	0.22 J	1 U	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	2000	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U
Vinyl chloride	1	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U
Xylene (total)	1000	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U	5 U	5 U	0.7 J	5 U	1.0 J
Total VOCs	-	ND	ND	ND	ND	0.2 J	4.7 J	3.3 J	ND	ND	ND	0.22 J	1.5	ND	ND	1.5 J	ND	1.7 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

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Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S	MW-13S	MW-14S	MW-14S	MW-14S	MW-14S	MW-14S	MW-14S	MW-15D
Sample Date	GROUNDWATER QUALITY	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/2/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	7/6/2007	6/29/2010	12/16/2010	12/29/2011	8/16/2012	12/20/2012	12/20/2004
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	33	5 U	NA
Benzene	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	0.23 J	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	0.088 J	1.3
Bromodichloromethane	1	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	NA	NA	4 U
Bromomethane	10	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	NA	NA	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	3.3 J	5 U	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA
Carbon tetrachloride	1	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	0.5	0.44 J	1 U	0.43 J	0.53 J	0.3 J	2 U
Chlorobenzene	50	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	NA	NA	5 U
Chloroethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	5 U
2-Chloroethyl vinyl ether	-	5 U	0.2 U	5 U	5 U	5 U	NA	NA	NA	NA	NA	0.2 U	NA	NA	NA	NA	NA	5 U
Chloroform	70	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.6	0.55 J	0.49 J	0.47 J	0.66 J	0.54 J	5 U
Chloromethane	-	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	NA	NA	5 U
Dibromochloromethane	1	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.3 U	1 U	1 U	1 U	1 U	1 U	5 U
1,1-Dichloroethane	50	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.3 U	1 U	1 U	1 U	1 U	1 U	5 U
1,2-Dichloroethane	2	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	0.3 U	1 U	1 U	1 U	0.19 J	1 U	0.6 J
1,1-Dichloroethene	1	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	5.1
cis-1,2-Dichloroethene	70	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.3 U	1 U	1 U	1 U	1 U	1 U	5 U
trans-1,2-Dichloroethene	100	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	5 U
1,2-Dichloropropane	1	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	NA	NA	1 U
cis-1,3-Dichloropropene	-	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.1 U	1 U	1 U	1 U	NA	NA	5 U
trans-1,3-Dichloropropene	-	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	NA	NA	5 U
Ethylbenzene	700	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	4 U
Methyl tert-butyl ether (MTBE)	70	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.3 U	1 U	1 U	1 U	NA	1 U	5 U
Methylene chloride	3	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	6.5 U	20 U	20 U	20 U	NA	NA	100 U
1,1,2,2-Tetrachloroethane	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	0.44 J	1 U	1 U
Tetrachloroethene	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	0.22 J	1 U
Toluene	600	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.3 U	1 U	1 U	0.19 J	1 U	1 U	2 J
1,1,1-Trichloroethane	30	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	5 U
1,1,2-Trichloroethane	3	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	3 U
Trichloroethene	1	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	0.8 J
Trichlorofluoromethane	2000	5 U	0.4 U	5 U	5 U	5 U	NA	NA	NA	NA	1 U	0.4 U	NA	NA	1 U	NA	NA	5 U
Vinyl chloride	1	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	2.5 J
Xylene (total)	1000	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	3 U	3 U	5 U
Total VOCs		ND	ND	ND	ND	ND	ND	ND	0.23 J	ND	ND	1.1	0.99 J	0.49 J	1.09 J	38.12 J	1.15 J	12.3 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID		MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-15D	MW-19S	MW-19S
Sample Date	GROUNDWATER QUALITY CRITERIA (7/22/2010) ug/L	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	7/11/2012	12/20/2012	6/29/2004	12/20/2004
(VOCs)		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Acetone	6000	NA	NA NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA NA
Benzene	1	4.3	0.9 J	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.16 J	1 U	0.7 J
Bromodichloromethane	1	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	10	1 U	10	1 U	1 U	1 U
Bromoform	4	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA.	NA NA	4 U	4 U
Bromomethane	10	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA.	NA NA	5 U	5 U
2- Butanone	300	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA	5 U	5 U	NA NA	NA NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA
Carbon tetrachloride	1	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U
Chlorobenzene	50	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U
Chloroethane	-	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	NA	NA	NA	5 U	5 U
Chloroform	70	5 U	5 U	5 U	5 U	0.3	0.3 J	0.4 J	0.4 J	0.41 J	0.33 J	0.42 J	0.47 J	0.29 J	0.25 J	0.27 J	5 U	5 U
Chloromethane	-	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U
Dibromochloromethane	1	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U
1,1-Dichloroethane	50	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2.6 J	2.8 J
1,2-Dichloroethane	2	1.1 J	2 U	2 U	0.5 J	0.3	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	0.22 J	2 U	2 U
1,1-Dichloroethene	1	7.6	4.1	3.0	2.2	2.2	1.2 J	1.1 J	1.1 J	1.8	2.2	1.1	1.4	0.88 J	1.4	2.9	2 U	2 U
cis-1,2-Dichloroethene	70	0.9 J	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1.7	1 U	1 U	1 U	1 U	1 U	5 U	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U
Ethylbenzene	700	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1.2 J	1 J
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U
Methylene chloride	3	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	NA	NA	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	5 U	5 U	0.6 J	5 U	0.3 U	5 U	5 U	5 U	1 U	0.41 J	1 U	1 U	1 U	1 U	1 U	5 U	1.9 J
1,1,1-Trichloroethane	30	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U
Trichloroethene	1	1.6	0.7 J	0.6 J	0.5 J	0.6	1 U	1 U	1 U	0.29 J	0.69 J	0.25 J	1 U	1 U	0.24 J	0.59 J	20	20
Trichlorofluoromethane	2000	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	0.21 J	NA	NA	5 U	5 U
Vinyl chloride	1	7.0	0.5 J	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U
Xylene (total)	1000	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U	2.9 J	2.5 J
Total VOCs		22.5 J	6.2 J	4.2 J	3.2 J	3.4	1.5 J	1.5 J	1.5 J	2.5 J	5.3 J	1.77 J	1.87 J	1.38 J	1.89 J	4.14 J	26.7 J	28.9 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

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regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-19S	MW-20S	MW-20S
Sample Date	GROUNDWATER QUALITY	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/10/2012	12/20/2012	6/28/2005	12/21/2005
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA
Benzene	1	1 U	0.5 J	1 U	1 U	0.2 U	1 U	1 U	0.8 J	1 U	1 U	1 U	1 U	0.37 J	0.13 J	0.088 J	1 U	1 U [1 U]
Bromodichloromethane	1	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]
Bromoform	4	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA	4 U	4 U [4 U]
Bromomethane	10	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U [5 U]
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA
Carbon tetrachloride	1	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U [2 U]
Chlorobenzene	50	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U [5 U]
Chloroethane	-	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	0.99 J	1 U	1 U	0.75 J	5 U	5 U [5 U]
2-Chloroethyl vinyl ether	-	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	NA	NA	NA	5 U	5 U [5 U]
Chloroform	70	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	0.3 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U [5 U]
Chloromethane	-	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U [5 U]
Dibromochloromethane	1	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U [5 U]
1,1-Dichloroethane	50	3.5 J	5 U	2.9 J	1.1 J	2.9	5 U	1.5 J	5 U	1.1	1 U	1.1	0.34 J	0.86 J	0.66 J	1 U	5 U	5 U [5 U]
1,2-Dichloroethane	2	2 U	2 U	2 U	2 U	0.2 U	2 U	2 U	0.3 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U [2 U]
1,1-Dichloroethene	1	2 U	2.6	2 U	2 U	0.5 U	2 U	2 U	2.0	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U [2 U]
cis-1,2-Dichloroethene	70	1.2 J	5 U	39	1.0 J	6.0	31	2.6 J	0.5 J	3.1	0.23 J	2.8	16	2.7	1.6	9	5 U	5 U [5 U]
trans-1,2-Dichloroethene	100	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U [5 U]
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U [1 U]
cis-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U [5 U]
trans-1,3-Dichloropropene	-	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U [5 U]
Ethylbenzene	700	0.4 J	4 U	2.0 J	4 U	0.4 U	0.7 J	0.5 J	4 U	1.0	1 U	0.77 J	2.0	1.6	0.97 J	2.7	4 U	4 U [4 U]
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U [5 U]
Methylene chloride	3	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U [3 U]
t-Butyl Alcohol (TBA)	100	100 U	9.3 J	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	20 U	20 U	NA	NA	100 U	100 U [100 U]
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]
Tetrachloroethene	1	0.5 J	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	0.26 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]
Toluene	600	5 U	5 U	0.7 J	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	0.18 J	1 U	1 U	1 U	5 U	5 U [5 U]
1,1,1-Trichloroethane	30	5 U	5 U	5 U	0.7 J	0.4 U	5 U	5 U	1.0 J	1 U	0.85 J	1 U	1 U	1 U	1 U	1 U	5 U	5 U [5 U]
1,1,2-Trichloroethane	3	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U [3 U]
Trichloroethene	1	15	3.3	1 U	4.0	3.3	1.1	3.4	3.8	2.1	1.7	0.94 J	0.73 J	1 U	0.64 J	0.12 J	1 U	1 U [1 U]
Trichlorofluoromethane	2000	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	0.68 J	NA	NA	5 U	5 U [5 U]
Vinyl chloride	1	5 U	0.4 J	5 U	5 U	0.2 U	5 U	5 U	0.8 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U [5 U]
Xylene (total)	1000	0.5 J	5 U	5.5	5 U	0.4	1.6 J	0.8 J	5 U	2.4 J	3 U	2.0 J	6.4	4.0	1.9 J	2.3 J	5 U	5 U [5 U]
Total VOCs	-	21.1 J	16.1 J	50.1 J	6.8	12.6	34.4 J	8.8 J	9.5 J	9.7 J	3 J	7.61 J	26.6 J	10.2 J	5.9 J	14.96 J	ND	ND [ND]

### Notes:

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Sample ID	NJ CLASS IIA	MW-20S	MW-20S	MW-20S	MW-20S	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D	MW-20D
Sample Date		6/21/2006	12/20/2006	7/6/2007	6/24/2008	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)			<u> </u>					<u> </u>						<u> </u>				
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	0.2 U	1 U	1 U	1.0	0.8 J	0.8 J	1 U	1.0	0.4 J	0.3 J	1 U [1 U]	1 U	0.35 J	0.32 J	0.26 J
Bromodichloromethane	1	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
Bromoform	4	4 U	4 U	0.2 U	4 U	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U [4 U]	1 U	1 U	1 U	1 U
Bromomethane	10	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	2 U	0.3 U	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U [2 U]	1 U	1 U	1 U	1 U
Chlorobenzene	50	5 U	5 U	0.2 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Chloroethane	-	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	-	5 U	5 U	0.2 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Chloroform	70	5 U	5 U	0.2 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Chloromethane	-	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Dibromochloromethane	1	5 U	5 U	0.3 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
1,1-Dichloroethane	50	5 U	5 U	0.3 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
1,2-Dichloroethane	2	2 U	2 U	0.3 U	2 U	2 U	2 U	2 U	2 U	2 U	0.3	2 U	2 U	2 U [2 U]	1 U	1 U	1 U	1 U
1,1-Dichloroethene	1	2 U	2 U	0.5 U	2 U	1.3 J	1.4 J	1.9 J	1.5 J	1.8 J	1.7	0.7 J	1.0	0.5 J [0.7 J]	0.85 J	0.25 J	0.82 J	1 U
cis-1,2-Dichloroethene	70	5 U	5 U	0.3 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
trans-1,2-Dichloroethene	100	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	0.1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
trans-1,3-Dichloropropene	-	5 U	5 U	0.2 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Ethylbenzene	700	4 U	4 U	0.4 U	4 U	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U [4 U]	1 U	1 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	0.6 J	5 U	0.3 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Methylene chloride	3	3 U	3 U	0.4 U	3 U	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U [3 U]	1 U	1 U	1 U	1 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	6.5 U	100 U	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U [100 U]	20 U	20 U	20 U	20 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U
Toluene	600	5 U	5 U	0.3 U	5 U	2.5 J	0.7 J	5 U	0.7 J	5 U	0.3 U	5 U	5 U	5 U [5 U]	1 U	1.1	1 U	1 U
1,1,1-Trichloroethane	30	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	3	3 U	3 U	0.2 U	3 U	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U [3 U]	1 U	1 U	1 U	1 U
Trichloroethene	1	1 U	1 U	0.4 U	1 U	1 U	1 U	0.5 J	1 U	0.5 J	0.4 U	1 U	1 U	1 U [1 U]	0.24 J	1 U	1 U	1 U
Trichlorofluoromethane	2000	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	1 U	1 U	1 U	1 U
Vinyl chloride	1	5 U	5 U	0.2 U	5 U	5 U	1.4 J	1.7 J	1.4 J	5 U	1.7	5 U	0.8 J	5 U [5 U]	1 U	1 U	0.52 J	1 U
Xylene (total)	1000	5 U	5 U	0.4 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U [5 U]	3 U	3 U	3 U	3 U
Total VOCs		0.6 J	ND	ND	ND	3.8 J	4.5 J	4.9 J	4.4 J	2.3 J	5.0	1.1 J	2.1	0.5 J [0.7 J]	1.1 J	1.7 J	1.66 J	0.26 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-20D	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-21S	MW-22D	MW-22D	MW-22D
Sample Date		12/29/2011	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/1/2009	12/23/2009	12/29/2011	12/20/2004	6/28/2005	12/21/2005
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)										<u>.</u>								
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	0.43 J	1 U	1 U	1 U	1 U	1 U	0.6 J	0.2 U	1 U	1 U	1 U	1 U	0.21 J	0.33 J	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	1 U	4 U	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	4 U	4 U	4 U
Bromomethane	10	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	2 U	2 U	2 U
Chlorobenzene	50	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Chloroethane	-	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
2-Chloroethyl vinyl ether	-	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Chloroform	70	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Chloromethane	-	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Dibromochloromethane	1	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
1,1-Dichloroethane	50	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
1,2-Dichloroethane	2	1 U	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	2 U	2 U	2 U
1,1-Dichloroethene	1	0.89 J	2 U	2 U	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	2 U	2 U	2 U
cis-1,2-Dichloroethene	70	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
trans-1,2-Dichloroethene	100	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
trans-1,3-Dichloropropene	-	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Ethylbenzene	700	1 U	4 U	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	4 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Methylene chloride	3	1 U	3 U	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U	3 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	20 U	100 U	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U	20 U	100 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	0.3 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	0.13 J	1 U	1 J	5 U	5 U
1,1,1-Trichloroethane	30	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
1,1,2-Trichloroethane	3	1 U	3 U	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	3 U	3 U	3 U
Trichloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 J	1 U
Trichlorofluoromethane	2000	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Vinyl chloride	1	0.64 J	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	5 U	5 U	5 U
Xylene (total)	1000	3 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	5 U	5 U	5 U
Total VOCs	-	1.96 J	ND	ND	ND	ND	ND	0.6 J	ND	0.3 J	ND	ND	ND	0.34 J	0.33 J	1 J	0.4 J	ND

### Notes:

U Not Detected Above Detection Limits
-- Not Sampled
Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22D	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I
Sample Date		6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)												<u> </u>						
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	0.7 J	0.2 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	0.22 J [0.27 J]	1 U	0.24 J	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	0.2 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	4 U	0.2 U	4 U	4 U [4 U]	4 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	4 U	4 U	4 U	4 U	4 U	4 U
Bromomethane	10	5 U	5 U	0.4 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	2 U	0.3 U	2 U	2 U [2 U]	2 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	2 U	2 U	2 U	2 U	2 U	2 U
Chlorobenzene	50	5 U	5 U	0.2 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Chloroethane	-	5 U	5 U	0.4 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
2-Chloroethyl vinyl ether	-	5 U	5 U	0.2 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Chloroform	70	5 U	5 U	0.2	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Chloromethane	-	5 U	5 U	0.4 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Dibromochloromethane	1	5 U	5 U	0.3 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
1,1-Dichloroethane	50	5 U	5 U	0.3 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloroethane	2	2 U	2 U	0.3 U	2 U	2 U [2 U]	2 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	2 U	2 U	2 U	2 U	2 U	2 U
1,1-Dichloroethene	1	2 U	2 U	1.2	2 U	0.6 J [0.5 J]	2 U	1 U [1 U]	1 U [1 U]	0.2 J [1 U]	1 U	1 U	2 U	2 U	2 U	2 U	2 U	2 U
cis-1,2-Dichloroethene	70	5 U	5 U	0.3 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
trans-1,2-Dichloroethene	100	5 U	5 U	0.4 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
1,2-Dichloropropane	1	1 U	1 U	0.5 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	0.1 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
trans-1,3-Dichloropropene	-	5 U	5 U	0.2 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Ethylbenzene	700	4 U	4 U	0.4 U	4 U	4 U [4 U]	4 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	4 U	4 U	4 U	4 U	4 U	4 U
Methyl tert-butyl ether (MTBE)	70	1.2 J	5 U	0.3 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	0.6 J	5 U
Methylene chloride	3	3 U	3 U	0.4 U	3 U	3 U [3 U]	3 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	1.4 J	3 U	3 U	3 U	3 U	3 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	6.5 U	100 U	100 U [100 U]	100 U	20 U [20 U]	20 U [20 U]	20 U [20 U]	20 U	3.0 J	100 U	100 U	100 U	100 U	100 U	100 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	0.4 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	0.4 U	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	1.4 J	5 U	0.3 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	1.4 J	1.4 J	5 U	5 U	5 U
1,1,1-Trichloroethane	30	5 U	5 U	0.4 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
1,1,2-Trichloroethane	3	3 U	3 U	0.2 U	3 U	3 U [3 U]	3 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	3 U	3 U	3 U	3 U	3 U	3 U
Trichloroethene	1	1 U	1 U	0.5	1 U	1 U [1 U]	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	0.6 J	1 U	1 U	1 U	1 U	1 U
Trichlorofluoromethane	2000	5 U	5 U	0.4 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Vinyl chloride	1	5 U	5 U	0.2 U	5 U	5 U [5 U]	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U
Xylene (total)	1000	1.2 J	5 U	0.4 U	5 U	5 U [5 U]	5 U	3 U [3 U]	3 U [3 U]	3 U [3 U]	3 U	3 U	5 U	5 U	0.6 J	5 U	5 U	5 U
Total VOCs		3.8 J	0.7 J	1.9	ND	0.6 J [0.5 J]	ND	ND [ND]	ND [ND]	0.42 J [0.27 J]	ND	3.24 J	2 J	1.4 J	2 J	ND	0.6 J	ND

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds

regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I	MW-22I	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S
Sample Date	GROUNDWATER QUALITY	7/6/2007	1/22/2008	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)							<u> </u>					<u> </u>						
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	0.2 U	1 U	1 U [1 U]	1 U	0.96 J	0.19 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	0.7 J [0.6 J]	0.2 U [0.2 U]	1 U [1 U]
Bromodichloromethane	1	0.2 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.2 U [0.2 U]	1 U [1 U]
Bromoform	4	0.2 U	4 U	4 U [4 U]	4 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U [4 U]	4 U [4 U]	0.2 U [0.2 U]	4 U [4 U]
Bromomethane	10	0.4 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	0.3 U	2 U	2 U [2 U]	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U [2 U]	2 U [2 U]	0.3 U [0.3 U]	2 U [2 U]
Chlorobenzene	50	0.2 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]
Chloroethane	-	0.4 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
2-Chloroethyl vinyl ether	-	0.2 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]
Chloroform	70	0.2 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [0.3 J]
Chloromethane	-	0.4 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
Dibromochloromethane	1	0.3 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]
1,1-Dichloroethane	50	0.3 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]
1,2-Dichloroethane	2	0.3 U	2 U	2 U [2 U]	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U [2 U]	2 U [2 U]	0.3 U [0.3 U]	2 U [2 U]
1,1-Dichloroethene	1	0.5 U	2 U	2 U [2 U]	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	0.7 J [2 U]	2 U [2 U]	0.5 U [0.5 U]	2 U [1.2 J]
cis-1,2-Dichloroethene	70	0.3 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]
trans-1,2-Dichloroethene	100	0.4 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
1,2-Dichloropropane	1	0.5 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.5 U [0.5 U]	1 U [1 U]
cis-1,3-Dichloropropene	-	0.1 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.1 U [0.1 U]	5 U [5 U]
trans-1,3-Dichloropropene	-	0.2 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]
Ethylbenzene	700	0.4 U	4 U	4 U [4 U]	4 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U [4 U]	4 U [4 U]	0.4 U [0.4 U]	4 U [4 U]
Methyl tert-butyl ether (MTBE)	70	0.3 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	0.8 J	2.0 J	0.8 J	5 U [0.6 J]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]
Methylene chloride	3	0.4 U	3 U	3 U [3 U]	3 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U [3 U]	3 U [3 U]	0.4 U [0.4 U]	3 U [0.3 J]
t-Butyl Alcohol (TBA)	100	6.5 U	100 U	100 U [100 U]	100 U	20 U	20 U	20 U	20 U	10 J	100 U	100 U	100 U	100 U	100 U [100 U]	100 U [100 U]	6.5 U [6.5 U]	100 U [100 U]
1,1,2,2-Tetrachloroethane	1	0.4 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.4 U [0.4 U]	1 U [1 U]
Tetrachloroethene	1	0.4 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.4 U [0.4 U]	1 U [1 U]
Toluene	600	0.3 U	5 U	5 U [5 U]	5 U	1 U	1.2	1 U	1 U	1 U	5 U	0.7 J	5 U	5 U	0.8 J [5 U]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]
1,1,1-Trichloroethane	30	0.4 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
1,1,2-Trichloroethane	3	0.2 U	3 U	3 U [3 U]	3 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U [3 U]	3 U [3 U]	0.2 U [0.2 U]	3 U [3 U]
Trichloroethene	1	0.4 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	0.4 U [0.4 U]	1 U [1 U]
Trichlorofluoromethane	2000	0.4 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
Vinyl chloride	1	0.2 U	5 U	5 U [5 U]	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]
Xylene (total)	1000	0.4 U	5 U	5 U [5 U]	5 U	3 U	3 U	3 U	3 U	3 U	5 U	5 U	5 U	5 U	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]
Total VOCs	-	ND	ND	ND [ND]	ND	0.96 J	1.4 J	ND	ND	10.0 J	ND	1.5 J	2.0 J	0.8 J	1.5 J [0.6 J]	0.7 J [0.6 J]	ND	ND [1.8 J]

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S	MW-22S	MW-23D	MW-23D	MW-23D	MW-23D	MW-23D	MW-23D	MW-23D	MW-23D	MW-23D	MW-23D
Sample Date	GROUNDWATER QUALITY	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	12/21/2004	6/28/2005	6/21/2006	12/20/2006	7/6/2006	12/27/2007	6/24/2008	12/19/2008	7/2/2009	12/23/2009
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	1 U	0.40 J	1 U	1 U	1.3	1 U	1 U	1 U	1.6	0.2 U	0.3 J	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	4 U	4 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U
Bromomethane	10	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U
2- Butanone	300	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	2 U	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U
Chlorobenzene	50	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U
Chloroethane	-	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U
2-Chloroethyl vinyl ether	-	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U
Chloroform	70	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.7 J	5 U	5 U	5 U	0.3	5 U	5 U	5 U	1 U	0.21 J
Chloromethane	-	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U
Dibromochloromethane	1	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U
1,1-Dichloroethane	50	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U
1,2-Dichloroethane	2	2 U	2 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U
1,1-Dichloroethene	1	2 U	2 U	1 U	1 U	1 U	1 U	1 U	2.0	2 U	3.5	2.3	1.2	0.7 J	0.9 J	2 U	1 U	1.1
cis-1,2-Dichloroethene	70	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U
trans-1,2-Dichloroethene	100	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U
1,2-Dichloropropane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U
trans-1,3-Dichloropropene	-	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U
Ethylbenzene	700	4 U	4 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	0.5 J	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U
Methylene chloride	3	3 U	3 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U
t-Butyl Alcohol (TBA)	100	100 U	100 U	20 U	20 U	20 U	20 U	20 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	5 U	5 U	1 U	1 U	1 U	1 U	1 U	0.9 J	0.4 J	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	0.64 J
1,1,1-Trichloroethane	30	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U
1,1,2-Trichloroethane	3	3 U	3 U	1 U	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U
Trichloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.6 J	0.7 J	1 U	0.4 U	1 U	1 U	1 U	1 U	0.44 J
Trichlorofluoromethane	2000	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U
Vinyl chloride	1	5 U	5 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	5 U	1.0 J	0.2 U	5 U	5 U	5 U	1 U	1 U
Xylene (total)	1000	5 U	5 U	3 U	3 U	3 U	3 U	3 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U
Total VOCs	-	ND	ND	ND	0.4 J	ND	ND	1.3	3.6 J	1 J	4.7 J	4.9 J	1.5	1.0 J	0.9 J	ND	ND	2.4 J

### Notes:

U Not Detected Above Detection Limits
-- Not Sampled
Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-23D	MW-23D	MW-23D	MW-23D	MW-23I	MW-23I	MW-231	MW-23I	MW-23I	MW-23I	MW-23I	MW-23I	MW-23I	MW-23I	MW-23I	MW-23I	MW-23I
Sample Date	GROUNDWATER QUALITY	6/30/2010	12/16/2010	12/29/2011	12/20/2012	6/29/2004	12/21/2004	6/28/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/2/2009	12/23/2009	6/30/2010	12/16/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	NA	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	0.15 J	0.36 J	1 U	1 U	1 U [1 U]	1 U	1.0	0.2 U	1 U	1 U	1 U	1 U [1 U]	0.39 J [0.43 J]	1 U [1 U]	1 U [1 U]
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Bromoform	4	1 U	1 U	1 U	NA	4 U	4 U	4 U [4 U]	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Bromomethane	10	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
2- Butanone	300	NA	NA	NA	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U	1 U	1 U	1 U	2 U	2 U	2 U [2 U]	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Chlorobenzene	50	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Chloroethane	-	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
2-Chloroethyl vinyl ether	-	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Chloroform	70	1 U	1 U	1 U	1 U	4.8	0.6 J	5 U [5 U]	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Chloromethane	-	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Dibromochloromethane	1	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
1,1-Dichloroethane	50	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
1,2-Dichloroethane	2	1 U	1 U	1 U	1 U	2 U	2.6	1.2 J [1.2 J]	0.6 J	2 U	0.3 U	2 U	0.8 J	2 U	1 U [1 U]	1 U [1 U]	0.26 J [0.30 J]	1 U [1 U]
1,1-Dichloroethene	1	0.40 J	1 U	0.49 J	1 U	0.7 J	1.3 J	1.0 J [1.0 J]	1.1 J	2 U	0.5 U	2 U	0.5 J	2 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
cis-1,2-Dichloroethene	70	1 U	1 U	1 U	1 U	0.9 J	5 U	5 U [5 U]	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
trans-1,2-Dichloroethene	100	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
1,2-Dichloropropane	1	1 U	1 U	1 U	NA	1 U	1 U	1 U [1 U]	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
cis-1,3-Dichloropropene	-	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
trans-1,3-Dichloropropene	-	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Ethylbenzene	700	1 U	1 U	1 U	1 U	4 U	4 U	4 U [4 U]	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Methylene chloride	3	1 U	1 U	1 U	1 U	2.4 J	3 U	3 U [3 U]	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
t-Butyl Alcohol (TBA)	100	20 U	20 U	20 U	NA	100 U	100 U	100 U [100 U]	100 U	100 U	6.5 U	100 U	100 U	14 J	20 U [20 U]	20 U [20 U]	20 U [20 U]	20 U [20 U]
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Toluene	600	1 U	1 U	1 U	1 U	5 U	2.5 J	0.5 J [0.5 J]	1.2 J	5 U	0.3 U	5 U	5 U	5 U	1 U [1 U]	0.23 J [0.35 J]	1 U [1 U]	1 U [1 U]
1,1,1-Trichloroethane	30	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
1,1,2-Trichloroethane	3	1 U	1 U	1 U	1 U	3 U	3 U	3 U [3 U]	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Trichloroethene	1	1 U	1 U	1 U	1 U	1.2	0.9 J	0.9 J [1.0 J]	0.7 J	1 U	0.4 U	1 U	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Trichlorofluoromethane	2000	1 U	1 U	1 U	NA	5 U	5 U	5 U [5 U]	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Vinyl chloride	1	1 U	1 U	1 U	1 U	5 U	5 U	5 U [5 U]	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]
Xylene (total)	1000	3 U	3 U	3 U	3 U	5 U	5 U	5 U [5 U]	0.9 J	5 U	0.4 U	5 U	5 U	5 U	3 U [3 U]	3 U [3 U]	3 U [3 U]	3 U [3 U]
Total VOCs	-	0.40 J	ND	0.64 J	0.36 J	10 J	7.9 J	3.6 J [3.7 J]	4.5 J	1.0	ND	ND	1.3 J	14 J	ND [ND]	0.62 J [0.78 J]	0.26 J [0.30 J]	ND [ND]

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-23I	MW-23I	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S	MW-23S
Sample Date	GROUNDWATER QUALITY	12/29/2011	12/20/2012	6/29/2004	12/21/2004	6/28/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/2/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	8/16/2012
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	33	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	48
Benzene	1	1 U	1 U	10 U [10 U]	25 U [25 U]	5 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	0.16 J	0.20 J	0.56 J
Bromodichloromethane	1	1 U	1 U	10 U [10 U]	25 U [25 U]	5 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	1 U	NA	40 U [40 U]	100 U [100 U]	20 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA
Bromomethane	10	1 U	NA	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA
2- Butanone	300	NA	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	2.8 J
Carbon Disulfide	700	NA	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	1 U
Carbon tetrachloride	1	1 U	1 U	20 U [20 U]	50 U [50 U]	10 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U
Chlorobenzene	50	1 U	NA	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA
Chloroethane	-	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	-	1 U	NA	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.2 U	5 U	5 U	5 U	NA	NA	NA	NA	NA	NA
Chloroform	70	1 U	1 U	29 J [23 J]	37 J [38 J]	32	22	14	14	6.1	3.6 J	0.8 J	1.0	1.6	4.0	4.5	3.8	1 U
Chloromethane	-	1 U	NA	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA
Dibromochloromethane	1	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	50	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	0.17 J	1 U	1 U
1,2-Dichloroethane	2	0.56 J	1 U	230 [220]	220 [220]	240	110	51	46	19	11	4.2	5.2	8.9	21	30	27	1 U
1,1-Dichloroethene	1	0.20 J	1 U	20 U [20 U]	50 U [50 U]	10 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	0.53 J
cis-1,2-Dichloroethene	70	1 U	1 U	68 [74]	50 J [49 J]	45	28	11	14	4.7 J	2.0 J	0.4 J	2.2	2.9	12	15	9.9	1 U
trans-1,2-Dichloroethene	100	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	0.6 J	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	0.20 J	1 U	1 U	1 U
1,2-Dichloropropane	1	1 U	NA	10 U [10 U]	25 U [25 U]	5 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA
cis-1,3-Dichloropropene	-	1 U	NA	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA
trans-1,3-Dichloropropene	-	1 U	NA	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA
Ethylbenzene	700	1 U	1 U	40 U [40 U]	100 U [100 U]	20 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	0.5 J	5 U	0.3 U	5 U	5 U	5 U	1 U	NA	1 U	1 U	1 U	NA
Methylene chloride	3	1 U	1 U	1500 [1400]	3200 [3200]	550	150	72	17	8.2	2.3 J	1.4 J	1 U	1 U	4.9	12	7.0	1 U
t-Butyl Alcohol (TBA)	100	20 U	NA	U]	U]	500 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	NA	20 U	20 U	20 U	NA
1,1,2,2-Tetrachloroethane	1	1 U	1 U	10 U [10 U]	25 U [25 U]	5 U	1.3	1.0	1.0	0.4 J	1 U	1 U	1 U	0.21 J	0.53 J	0.51 J	0.58 J	1 U
Tetrachloroethene	1	1 U	1 U	10 U [10 U]	25 U [25 U]	5 U	0.8 J	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U
1,1,1-Trichloroethane	30	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	1.0 J	0.6 J	0.6	5 U	5 U	5 U	1 U	1 U	1 U	1 U	0.29 J	1 U
1,1,2-Trichloroethane	3	1 U	1 U	30 U [30 U]	75 U [75 U]	15 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1	1 U	1 U	37 [36]	22 J [21 J]	24	13	5.4	6.4	2.1	1.4	1 U	0.68 J	0.66 J	5.2	8.9	1 U	1 U
Trichlorofluoromethane	2000	1 U	NA	50 U [11 J]	120 U [120 U]	25 U	1.2 J	5 U	0.4 U	5 U	5 U	5 U	NA	NA	NA	NA	6.6	NA
Vinyl chloride	1	1 U	1 U	50 U [50 U]	120 U [120 U]	25 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U
Xylene (total)	1000	3 U	3 U	50 U [50 U]	120 U [120 U]	25 U	0.6 J	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U
Total VOCs	-	0.76 J	33	J]	J]	891	329 J	155 J	99	40.5 J	20.3	6.8 J	9.1 J	14.3 J	47.8 J	71.2 J	55.4 J	51.89 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

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regulatory criteria Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-23S	MW-24	MW-24	MW-24	MW-24	MW-24	MW-24	MW-24	MW-24	MW-24	MW-25	MW-25	MW-25	MW-25	MW-25	MW-25	MW-25
Sample Date	GROUNDWATER QUALITY	12/20/2012	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010	12/29/2011	7/10/2012	12/20/2012	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	5 U [5 U]	NA	NA	NA	NA	NA	NA	NA	5 U [5 U]	5 U	NA	NA	NA	NA	NA	NA	NA
Benzene	1	0.12 J [0.10 J]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.14 J [0.13 J]	0.5 J	1 U	1 U	1 U	0.41 J	1 U	0.44 J	1 U
Bromodichloromethane	1	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	NA	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	4 U	4 U	1 U	1 U	1 U	1 U	1 U
Bromomethane	10	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U
2- Butanone	300	5 U [5 U]	NA	NA	NA	NA	NA	NA	NA	5 U [5 U]	5 U	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	1 U [1 U]	NA	NA	NA	NA	NA	NA	NA	1 U [1 U]	1 U	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U [1 U]	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	0.5 J	0.7 J	1 U	0.21 J	0.55 J	1 U	0.40 J
Chlorobenzene	50	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U
Chloroethane	-	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	=	NA	5 U	5 U	NA	NA	NA	NA	1 U	NA [NA]	NA	5 U	5 U	NA	NA	NA	NA	NA
Chloroform	70	0.57 J [0.55 J]	5 U	5 U	1 U	0.25 J	1 U	1 U	1 U	0.13 J [0.13 J]	1 U	5.0	1.9 J	5.7	2.4	3.0	0.49 J	1.4
Chloromethane	=	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	50	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	0.41 J	1 U	0.24 J	1 U
1,2-Dichloroethane	2	5.5 [5.5]	2 U	2 U	1 U	1 U	1 U	1.4	1 U	1 U [1 U]	1 U	4.5	0.6 J	2.8	34	4.7	45	8.2
1,1-Dichloroethene	1	1 U [1 U]	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U
cis-1,2-Dichloroethene	70	1.1 [1.1]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	8.7	1.3 J	13	14	14	3.0	4.3
trans-1,2-Dichloroethene	100	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	0.28 J	0.23 J	0.27 J	1 U
1,2-Dichloropropane	1	NA	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	1 U	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	•	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U
trans-1,3-Dichloropropene	-	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA [NA]	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U
Ethylbenzene	700	1 U [1 U]	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	1 U [1 U]	5 U	5 U	1 U	NA	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	NA	1 U	1 U	1 U
Methylene chloride	3	1.8 [1.7]	3 U	3 U	1 U	1 U	1 U	1.7	1 U	1 U [1 U]	1 U	3 U	3 U	1 U	2.5	0.82 J	1.1	1.6
t-Butyl Alcohol (TBA)	100	NA	100 U	100 U	20 U	NA	20 U	20 U	20 U	NA [NA]	NA	100 U	100 U	20 U	NA	20 U	20 U	5.2 J
1,1,2,2-Tetrachloroethane	1	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	0.13 J	0.21 J [0.18 J]	1 U	1 U	1 U	1 U	0.34 J	1 U	1 U	0.11 J
Tetrachloroethene	1	1 U [1 U]	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	0.12 J	0.6 J	1 U	0.63 J	0.55 J	1.1	0.39 J	0.41 J
Toluene	600	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U
1,1,1-Trichloroethane	30	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	0.7 J	5 U	1 U	1.3	0.58 J	1.2	1 U
1,1,2-Trichloroethane	3	1 U [1 U]	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U
Trichloroethene	1	1.2 [1.2]	1 U	1 U	1 U	1 U	1 U	0.52 J	1 U	1 U [1 U]	1 U	8.0	1.2	9.7	41	9.6	37	1 U
Trichlorofluoromethane	2000	NA	5 U	5 U	NA	NA	NA	NA	1 U	NA [NA]	NA	5 U	5 U	NA	NA	NA	NA	4.8
Vinyl chloride	1	1 U [1 U]	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U
Xylene (total)	1000	3 U [3 U]	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U [3 U]	3 U	5 U	5 U	3 U	3 U	3 U	1.0 J	3 U
Total VOCs	-	J]	ND	ND	ND	0.25 J	ND	3.62 J	0.13 J	0.48 J [0.44 J]	0.62 J	28 J	5.7 J	31.8 J	97.4 J	34.6 J	90.1 J	26.4 J

### Notes:

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Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-25	MW-25	MW-26	MW-26	MW-26	MW-26	MW-26	MW-26	MW-26	MW-26	MW-26	MW-27	MW-27	MW-27	MW-27	MW-27	MW-27
Sample Date	GROUNDWATER QUALITY	7/11/2012	12/20/2012	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/11/2012	12/20/2012	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)												<u> </u>						
Acetone	6000	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	1 U	1 U	1 U	1 U	0.74 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	NA	NA	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA	4 U	4 U	1 U	1 U	1 U	1 U
Bromomethane	10	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U
2- Butanone	300	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	1 U	1 U	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	0.55 J	1 U	2 U	9.8	3.4	1.6	1 U	1.0	1.3	1.6	0.44 J	2 U	2 U	1 U	1 U	1 U	1 U
Chlorobenzene	50	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U
Chloroethane	-	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	-	NA	NA	5 U	5 U	NA	NA	NA	NA	1 U	NA	NA	5 U	5 U	NA	NA	NA	NA
Chloroform	70	1.4	0.56 J	2.8 J	10	8.0	3.8	1.2	0.67 J	2.6	3.9	0.75 J	5 U	0.7 J	2.0	2.1	0.48 J	0.20 J
Chloromethane	-	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	50	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U
1,2-Dichloroethane	2	6.2	7.6	20	2	40	16	120	28	6.8	11	10	2 U	2 U	1 U	1 U	1 U	1 U
1,1-Dichloroethene	1	1 U	1 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	2 U	1 U	1 U	1 U	1 U
cis-1,2-Dichloroethene	70	8.2	29	8.2	5 U	1.4	2.1	11	4.4	0.34 J	6	3.9	5 U	4.3 J	1 U	1 U	1 U	1 U
trans-1,2-Dichloroethene	100	1 U	1.5	5 U	5 U	1 U	1 U	0.22 J	1 U	1 U	0.44 J	1 U	5 U	0.5 J	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U
trans-1,3-Dichloropropene	-	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U
Ethylbenzene	700	1 U	1 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	4 U	4 U	1 U	1 U	1 U	1.0
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	5 U	5 U	1 U	NA	1 U	1 U	1 U	1 U	1 U	5 U	5 U	1 U	NA	1 U	1 U
Methylene chloride	3	1.9	2.6	1.2 J	3 U	1.1	0.85 J	16	2.1	0.37 J	1.1	1.3	3 U	3 U	1 U	1 U	1 U	1 U
t-Butyl Alcohol (TBA)	100	NA	NA	100 U	100 U	20 U	NA	20 U	20 U	20 U	NA	NA	100 U	100 U	20 U	NA	20 U	20 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	3.0	3.0	1.8	2.4	0.43 J	1.5	1.7	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	0.38 J	0.73 J	0.4 J	4.7	0.79 J	0.62 J	0.71 J	0.92 J	0.42 J	0.44 J	0.28 J	5.2	1 U	1 U	0.65 J	0.35 J	0.60 J
Toluene	600	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U
1,1,1-Trichloroethane	30	1 U	0.21 J	5 U	5 U	1 U	1 U	1 U	0.35 J	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	3	1 U	1 U	3 U	3 U	1 U	1 U	0.34 J	1 U	1 U	1 U	1 U	3 U	3 U	1 U	1 U	1 U	1 U
Trichloroethene	1	6	27	6.8	1.2	2.1	1.6	9.0	17	1 U	7.3	4.7	10	57	16	8.5	4.7	4.7
Trichlorofluoromethane	2000	NA	NA	5 U	5 U	NA	NA	NA	NA	1.2	NA	NA	5 U	5 U	NA	NA	NA	NA
Vinyl chloride	1	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U
Xylene (total)	1000	3 U	3 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U	5 U	5 U	3 U	3 U	3 U	2.4 J
Total VOCs	-	24.63 J	69.2 J	39.4 J	30.7	59.8 J	28.4 J	161.6 J	54.9 J	14.5 J	33.48 J	21.37 J	15.2	62.5 J	18	11.3 J	5.53 J	8.9 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	MW-27	MW-27	MW-27	MW-28	MW-28	MW-28	MW-28	MW-28	MW-28	MW-28	MW-28	MW-28	PZ-1S	PZ-1S	PZ-1S	PZ-1S	PZ-1S
Sample Date	GROUNDWATER QUALITY	12/29/2011	7/10/2012	12/20/2012	6/24/2008	12/19/2008	6/30/2009	12/23/2009	6/29/2010	12/16/2010	12/29/2011	7/11/2012	12/20/2012	12/19/2008	7/1/2009	12/23/2009	6/30/2010	12/16/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)		<u> </u>		· 0,		- J.		· 0,		<u></u>	- J.		<i>.</i>	- 5,	<u> </u>		- <u>U</u>	J.
Acetone	6000	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Benzene	1	0.31 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 J	1 U	1 U	1 U	0.39 J
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	1 U	NA	NA	4 U	4 U	1 U	1 U	1 U	1 U	1 U	NA	NA	4 U	1 U	1 U	1 U	1 U
Bromomethane	10	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	1 U	1 U	1 U	1 U
2- Butanone	300	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	1 U	1 U	NA	NA	NA	NA	NA	NA	NA	1 U	1 U	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U	1 U	1 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	1 U	1 U	1 U	1 U
Chlorobenzene	50	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	1 U	1 U	1 U	1 U
Chloroethane	-	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	-	NA	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	5 U	1 U	NA	1 U	NA
Chloroform	70	0.27 J	1 U	1 U	5 U	5 U	1 U	0.67 J	1 U	0.59 J	1 U	1 U	1 U	3.0 J	0.46 J	0.16 J	1 U	0.31 J
Chloromethane	-	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	1 U	1 U	1 U	1 U
Dibromochloromethane	1	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U	1 U
1,1-Dichloroethane	50	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U	1 U
1,2-Dichloroethane	2	1 U	1 U	1 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	20	4.9	6.3	3.2	32
1,1-Dichloroethene	1	1 U	1 U	1 U	2 U	2 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	2 U	1 U	1 U	1 U	1 U
cis-1,2-Dichloroethene	70	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	7.1	0.94 J	0.39 J	0.61 J	0.99 J
trans-1,2-Dichloroethene	100	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U	1 U
1,2-Dichloropropane	1	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	1 U	1 U	1 U	1 U
trans-1,3-Dichloropropene	-	1 U	NA	NA	5 U	5 U	1 U	1 U	1 U	1 U	1 U	NA	NA	5 U	1 U	1 U	1 U	1 U
Ethylbenzene	700	1 U	1 U	1 U	4 U	4 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	4 U	1 U	1 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	1 U	5 U	5 U	1 U	NA	1 U	1 U	1 U	1 U	1 U	5 U	1 U	NA	1 U	1 U
Methylene chloride	3	1 U	1 U	1 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.9 J	1 U	2.4	1 U	0.69 J
t-Butyl Alcohol (TBA)	100	20 U	NA	NA	100 U	100 U	20 U	NA	20 U	20 U	20 U	NA	NA	100 U	20 U	NA	20 U	20 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U
Tetrachloroethene	1	0.27 J	0.26 J	0.47 J	0.7 J	2.0	1.0	0.27 J	0.39 J	1 U	0.50 J	0.57 J	0.67 J	1 U	1 U	1 U	1 U	0.28 J
Toluene	600	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U	1 U
1,1,1-Trichloroethane	30	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.7 J	1 U	1 U	1 U	0.80 J
1,1,2-Trichloroethane	3	1 U	1 U	1 U	3 U	3 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	3 U	1 U	1 U	1 U	1 U
Trichloroethene	1	1 U	13	9.5	1 U	1 U	0.59 J	3.2	0.29 J	4.6	1 U	0.64 J	0.64 J	17	2.1	1.8	2.3	11
Trichlorofluoromethane	2000	9.6	NA	NA	5 U	5 U	NA	NA	NA	NA	0.47 J	NA	NA	5 U	1 U	1 U	1 U	1 U
Vinyl chloride	1	1 U	1 U	1 U	5 U	5 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	5 U	1 U	1 U	1 U	1 U
Xylene (total)	1000	3 U	3 U	3 U	5 U	5 U	3 U	3 U	3 U	3 U	3 U	3 U	3 U	0.2 J	3 U	3 U	3 U	3 U
Total VOCs	-	10.5 J	13.26 J	9.97 J	0.7 J	2	1.6 J	4.1 J	0.68 J	5.19 J	0.97 J	1.21 J	1.31 J	49.3 J	8.4 J	11.1 J	6.1 J	46.7 J

### Notes:

U Not Detected Above Detection Limits

-- Not Sampled

Bolded value indicates a detect above detection limits Red bolded value indicates a detection that exceeds regulatory criteria

Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	PZ-1S	PZ-1S	PZ-1S	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1M
Sample Date	GROUNDWATER QUALITY	12/29/2011	7/10/2012	12/20/2012	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/1/2009	12/23/2009	6/30/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)															<u> </u>		
Acetone	6000	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	1 U	1 U [1 U]	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	0.2 U [0.2 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U
Bromodichloromethane	1	1 U	1 U	0.38 J	1 U [1 U]	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	0.2 U [0.2 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U
Bromoform	4	1 U	NA	NA	4 U [4 U]	4 U	4 U	4 U [4 U]	4 U [4 U]	4 U [4 U]	0.2 U [0.2 U]	4 U [4 U]	4 U [4 U]	4 U [4 U]	1 U [1 U]	1 U	1 U
Bromomethane	10	1 U	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
2- Butanone	300	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	1 U	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U	1 U	1 U	2 U [2 U]	2 U	2 U	2 U [2 U]	2 U [2 U]	2 U [2 U]	0.3 U [0.3 U]	2 U [2 U]	2 U [2 U]	2 U [2 U]	1 U [1 U]	1 U	1 U
Chlorobenzene	50	1 U	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
Chloroethane		1 U	1 U	1 U	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
2-Chloroethyl vinyl ether	-	1 U	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	NA
Chloroform	70	0.18 J	1 U	1.6	1.8 J [1.8 J]	2.4 J	1.8 J	2.0 J [2.1 J]	1.0 J [1.3 J]	1.1 J [1.0 J]	0.7 [0.7]	1.2 J [1.1 J]	0.6 J [0.6 J]	0.9 J [0.8 J]	0.68 J [0.58 J]	0.21 J	0.33 J
Chloromethane	-	1 U	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
Dibromochloromethane	1	1 U	1 U	1 U	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
1,1-Dichloroethane	50	1 U	1 U	1 U	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
1,2-Dichloroethane	2	5.0	0.24 J	4.8	14 [15]	34	11	44 [46]	12 [12]	35 [36]	5.6 [6.0]	55 [54]	8.9 [9.3]	90 [81]	29 [28]	19	7.5
1,1-Dichloroethene	1	1 U	1 U	1 U	0.6 J [0.5 J]	0.6 J	0.8 J	2 U [2 U]	1.0 J [0.9 J]	0.6 J [0.6 J]	1.1 [1.0]	2 U [2 U]	1.0 J [0.9 J]	2 U [2 U]	1 U [1 U]	0.31 J	0.39 J
cis-1,2-Dichloroethene	70	0.33 J	0.23 J	1.9	14 [15]	9	5.8	8.5 [8.5]	3.2 J [3.0 J]	6.3 [6.1 J]	4.1 [4.1]	11 [11]	1.0 J [1.0 J]	11 [10]	4.4 [4.2]	3.1	1.3
trans-1,2-Dichloroethene	100	1 U	1 U	1 U	0.3 J [0.3 J]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
1,2-Dichloropropane	1	1 U	NA	NA	1 U [1 U]	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	0.5 U [0.5 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U
cis-1,3-Dichloropropene	-	1 U	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.1 U [0.1 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
trans-1,3-Dichloropropene	-	1 U	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
Ethylbenzene	700	1 U	1 U	1 U	4 U [4 U]	4 U	4 U	4 U [4 U]	4 U [4 U]	4 U [4 U]	0.4 U [0.4 U]	4 U [4 U]	4 U [4 U]	4 U [4 U]	1 U [1 U]	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	1 U	5 U [5 U]	5 U	5 U	5 U [5 U]	0.6 J [0.6 J]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
Methylene chloride	3	0.72 J	1 U	1.6	1.7 J [1.8 J]	1 J	3 U	1.1 J [1.2 J]	0.8 J [0.7 J]	0.7 J [0.7 J]	0.8 [0.8]	0.7 J [0.7 J]	0.8 J [0.8 J]	0.9 J [0.9 J]	1 U [1 U]	1 U	0.72 J
t-Butyl Alcohol (TBA)	100	20 U	NA	NA	100 U [100 U]	100 U	100 U	100 U [100 U]	100 U [100 U]	100 U [100 U]	6.5 U [6.5 U]	100 U [100 U]	100 U [100 U]	7.9 J [10 J]	20 U [20 U]	20 U	20 U
1,1,2,2-Tetrachloroethane	1	1 U	0.53 J	1 U	1.5 [1.3]	1.1	0.9 J	0.8 J [0.8 J]	1 U [0.5 J]	0.6 J [0.5 J]	0.4 [0.4]	0.7 J [0.7 J]	1 U [1 U]	0.3 J [0.3 J]	1 U [0.29 J]	0.16 J	1 U
Tetrachloroethene	1	1 U	0.6 J	0.25 J	0.3 J [1 U]	1 U	1 U	1 U [1 U]	1 U [1 U]	1 U [1 U]	0.4 U [0.4 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U [1 U]	1 U	1 U
Toluene	600	1 U	1 U	1 U	5 U [5 U]	2.4 J	0.8 J	5 U [5 U]	0.4 J [0.4 J]	5 U [5 U]	0.3 U [0.3 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1.3	1 U
1,1,1-Trichloroethane	30	1 U	1 U	1 U	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
1,1,2-Trichloroethane	3	1 U	1 U	1 U	3 U [3 U]	3 U	3 U	3 U [3 U]	3 U [3 U]	3 U [3 U]	0.2 U [0.2 U]	3 U [3 U]	3 U [3 U]	3 U [3 U]	1 U [1 U]	1 U	1 U
Trichloroethene	1	1 U	0.68 J	4.2	8 [8.1]	4.1	4.2	4.4 [4.1]	2.6 [2.4]	3 [2.9]	2.7 [2.8]	6.4 [5.8]	1.2 [1.2]	5.5 [5.6]	2.7 [2.4]	1.6	0.7 J
Trichlorofluoromethane	2000	1.1	NA	NA	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
Vinyl chloride	1	1 U	1 U	1 U	5 U [5 U]	5 U	5 U	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.2 U [0.2 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	1 U [1 U]	1 U	1 U
Xylene (total)	1000	3 U	3 U	3 U	5 U [5 U]	5 U	0.6 J	5 U [5 U]	5 U [5 U]	5 U [5 U]	0.4 U [0.4 U]	5 U [5 U]	5 U [5 U]	5 U [5 U]	3 U [3 U]	3 U	3 U
Total VOCs	-	7.33 J	2.28 J	14.73 J	42.2 J [43.8 J]	54.6 J	25.9 J	60.8 J [62.7 J]	21.6 J [21.8 J]	47.3 J [47.8 J]	15.4 [15.8]	75 J [73.3 J]	13.5 [13.8]	116.5 J [108.6 J]	36.8 J [35.5 J]	25.7 J	10.9 J

U Not Detected Above Detection Limits

-- Not Sampled

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Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	WCC-1M	WCC-1M	WCC-1M	WCC-1M	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-1S
Sample Date	GROUNDWATER QUALITY	12/16/2010	12/29/2011	7/11/2012	12/20/2012	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	12/27/2007	6/24/2008	12/19/2008	7/2/2009	12/23/2009	6/30/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)			- U		<i></i>		- J.	- J.		<i></i>		<u> </u>	- J.		<u> </u>		<i></i>	- 0.
Acetone	6000	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	0.22 J	1 U	0.45 J	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Bromoform	4	1 U	1 U	NA	NA	4 U	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U [1 U]	1 U
Bromomethane	10	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
2- Butanone	300	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	1 U	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U [1 U]	1 U
Chlorobenzene	50	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Chloroethane	-	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
2-Chloroethyl vinyl ether	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Chloroform	70	0.54 J	0.41 J	0.32 J	0.54 J	17	5 U	5 U	5 U	5 U	5 U	1.2	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Chloromethane	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Dibromochloromethane	1	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
1,1-Dichloroethane	50	0.24 J	1 U	1 U	1 U	0.8 J	5 U	0.8 J	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
1,2-Dichloroethane	2	56	3.5	3.7	49	74	2 U	77	2 U	2 U	1.0 J	6.5	2 U	2 U	2 U	1 U	1 U [1 U]	0.51 J
1,1-Dichloroethene	1	1 U	0.48 J	0.47 J	1 U	2 U	2 U	2 U	2 U	2 U	2 U	0.5 U	2 U	2 U	2 U	1 U	1 U [1 U]	1 U
cis-1,2-Dichloroethene	70	9.0	1.4	1.5	11	26	5 U	17	5 U	5 U	5 U	1.6	5 U	5 U	5 U	1 U	1 U [1 U]	0.49 J
trans-1,2-Dichloroethene	100	1 U	1 U	1 U	1 U	0.8 J	5 U	0.7 J	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
1,2-Dichloropropane	1	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
cis-1,3-Dichloropropene	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
trans-1,3-Dichloropropene	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Ethylbenzene	700	1 U	1 U	1 U	1 U	4 U	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U [1 U]	1 U
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	1 U	1 U	0.6 J	5 U	0.9 J	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Methylene chloride	3	0.60 J	0.50 J	1 U	1 U	8.7	3 U	1.3 J	3 U	3 U	3 U	1.4	3 U	3 U	3 U	1 U	1 U [1 U]	1 U
t-Butyl Alcohol (TBA)	100	20 U	20 U	NA	NA	100 U	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	100 U	20 U	20 U [20 U]	20 U
1,1,2,2-Tetrachloroethane	1	0.30 J	0.19 J	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1.0	1 U	0.5 J	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U [1 U]	1 U
Toluene	600	1 U	1 U	1 U	1 U	5 U	5 U	0.5 J	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
1,1,1-Trichloroethane	30	1 U	1 U	1 U	1 U	0.9 J	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
1,1,2-Trichloroethane	3	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U [1 U]	1 U
Trichloroethene	1	4.7	1 U	2.1	4.8	14	1 U	7.7	1 U	1 U	1 U	2.3	1 U	1 U	1 U	1 U	1 U [1 U]	0.82 J
Trichlorofluoromethane	2000	1 U	1.6	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Vinyl chloride	1	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U [1 U]	1 U
Xylene (total)	1000	3 U	3 U	3 U	3 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	3 U [3 U]	3 U
Total VOCs		71.4 J	8.3 J	8.09 J	65.79 J	144 J	ND	106.4 J	ND	ND	1.0 J	13.0	ND	ND	ND	ND	ND [ND]	1.82 J

### Notes:

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-- Not Sampled

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Historic groundwater data are obtained from the 2012 Annual Groundwater Report (Arcadis, 2012)

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Sample ID	NJ CLASS IIA	WCC-1S	WCC-1S	WCC-1S	WCC-1S	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M	WCC-3M
Sample Date	GROUNDWATER QUALITY	12/16/2010	12/29/2011	7/11/2012	12/20/2012	6/29/2004	12/20/2004	6/28/2005	12/21/2005	6/21/2006	12/20/2006	7/6/2007	1/22/2008	6/24/2008	12/19/2008	7/2/2009	12/23/2009	6/30/2010
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)																		
Acetone	6000	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Benzene	1	1 U	1 U	0.13 J	0.17 J	0.5 J	1 U	1 U	1 U	1 U	1 U	0.3	0.4 J	0.3 J	1 U	1 U	0.45 J	0.32 J
Bromodichloromethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.2 U	1 U	1 U	1 U	1 U	1 U	1 U
Bromoform	4	1 U	1 U	NA	NA	4 U	4 U	4 U	4 U	4 U	4 U	0.2 U	4 U	4 U	4 U	1 U	1 U	1 U
Bromomethane	10	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U
2- Butanone	300	NA	NA	5 U	5 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon Disulfide	700	NA	NA	1 U	1 U	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
Carbon tetrachloride	1	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U
Chlorobenzene	50	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U
Chloroethane	-	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U
Chloroform	70	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	0.3 J	1 U	0.18 J	0.28 J
Chloromethane	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U
Dibromochloromethane	1	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U
1,1-Dichloroethane	50	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	5 U	1 U	1 U	1 U
1,2-Dichloroethane	2	1 U	23	5.3	0.92 J	2 U	2 U	2 U	2 U	2 U	2 U	0.3 U	2 U	2 U	2 U	1 U	1 U	1 U
1,1-Dichloroethene	1	1 U	1 U	1 U	1 U	2 U	2 U	2 U	2 U	2 U	2 U	0.6	0.6 J	0.5 J	2 U	1 U	1 U	1 U
cis-1,2-Dichloroethene	70	1 U	7.9	3.3	0.41 J	5 U	5 U	5 U	5 U	5 U	5 U	0.3 U	5 U	5 U	0.2 J	1 U	0.20 J	1 U
trans-1,2-Dichloroethene	100	1 U	0.22 J	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U
1,2-Dichloropropane	1	1 U	1 U	NA	NA	1 U	1 U	1 U	1 U	1 U	1 U	0.5 U	1 U	1 U	1 U	1 U	1 U	1 U
cis-1,3-Dichloropropene	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.1 U	5 U	5 U	5 U	1 U	1 U	1 U
trans-1,3-Dichloropropene	-	1 U	1 U	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U
Ethylbenzene	700	1 U	1 U	1 U	1 U	0.7 J	4 U	4 U	4 U	4 U	4 U	0.4 U	4 U	4 U	4 U	1 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	1.4 J	0.6 J	0.6	1.8 J	0.8 J	0.3 J	1 U	0.74 J	0.70 J
Methylene chloride	3	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U	3 U	0.4 U	3 U	3 U	3 U	1 U	1 U	1 U
t-Butyl Alcohol (TBA)	100	20 U	20 U	NA	NA	100 U	100 U	100 U	100 U	100 U	100 U	6.5 U	100 U	100 U	9.8 J	20 U	20 U	20 U
1,1,2,2-Tetrachloroethane	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.5 J	0.8	1.0	1.0 J	0.7 J	1.3	0.88 J	0.50 J
Tetrachloroethene	1	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	1 U	1 U
Toluene	600	1 U	1 U	1 U	1 U	5 U	1.5 J	5 U	5 U	5 U	5 U	0.5	5 U	5 U	5 U	1 U	0.22 J	1 U
1,1,1-Trichloroethane	30	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U
1,1,2-Trichloroethane	3	1 U	1 U	1 U	1 U	3 U	3 U	3 U	3 U	3 U	3 U	0.2 U	3 U	3 U	3 U	1 U	1 U	1 U
Trichloroethene	1	1 U	1 U	1.7	0.52 J	1 U	1 U	1 U	1 U	1 U	1 U	0.4 U	1 U	1 U	1 U	1 U	0.34 J	0.26 J
Trichlorofluoromethane	2000	1 U	3.4	NA	NA	5 U	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	1 U	1 U	1 U
Vinyl chloride	1	1 U	1 U	1 U	1 U	5 U	5 U	5 U	5 U	5 U	5 U	0.2 U	5 U	5 U	5 U	1 U	1 U	1 U
Xylene (total)	1000	3 U	3 U	3 U	3 U	5.7	5 U	5 U	5 U	5 U	5 U	0.4 U	5 U	5 U	5 U	3 U	0.44 J	3 U
Total VOCs		ND	35.5 J	10.43 J	2.02 J	6.9 J	1.5 J	ND	ND	1.4 J	1.1 J	2.8	3.8 J	2.6 J	11.3 J	1.3	3.5 J	2.06 J

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Sample ID Sample Date	NJ CLASS IIA GROUNDWATER QUALITY	WCC-3M 12/16/2010	WCC-3M 12/29/2011	WCC-3M 7/11/2012	WCC-3M 1/9/2013	IW-3S 7/12/2012	IW-4S 7/12/2012	IW1-BT-2 7/10/2012
Unit	CRITERIA (7/22/2010) ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
(VOCs)			-0/-	-87 -		6/ -	-6/ -	6/ -
Acetone	6000	NA	NA	5 U [5 U]	5 U	51	5 U	5 U
Benzene	1	NA	NA	0.3 J [0.25 J]	0.24 J	0.27 J	1 U	1 U
Bromodichloromethane	1	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
Bromoform	4	NA	NA	NA	NA	NA	NA	NA
Bromomethane	10	NA	NA	NA	NA	NA	NA	NA
2- Butanone	300	NA	NA	1 U [1 U]	1 U	2.8 J	5 U	5 U
Carbon Disulfide	700	NA	NA	1 U [1 U]	1 U	0.77 J	1 U	1 U
Carbon tetrachloride	1	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
Chlorobenzene	50	NA	NA	NA	NA	NA	NA	NA
Chloroethane	-	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
2-Chloroethyl vinyl ether	-	NA	NA	NA	NA	NA	NA	NA
Chloroform	70	NA	NA	0.25 J [0.31 J]	0.38 J	1.5	0.26 J	1 U
Chloromethane	=	NA	NA	NA	NA	NA	NA	NA
Dibromochloromethane	1	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
1,1-Dichloroethane	50	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
1,2-Dichloroethane	2	NA	NA	1 U [1 U]	1 U	140	1 U	1 U
1,1-Dichloroethene	1	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
cis-1,2-Dichloroethene	70	NA	NA	1 U [1 U]	1 U	0.61 J	0.37 J	38
trans-1,2-Dichloroethene	100	NA	NA	1 U [1 U]	1 U	1 U	1 U	0.69 J
1,2-Dichloropropane	1	NA	NA	NA	NA	NA	NA	NA
cis-1,3-Dichloropropene	-	NA	NA	NA	NA	NA	NA	NA
trans-1,3-Dichloropropene	=	NA	NA	NA	NA	NA	NA	NA
Ethylbenzene	700	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
Methyl tert-butyl ether (MTBE)	70	NA	NA	0.22 J [0.28 J]	1 U	1 U	1 U	1 U
Methylene chloride	3	NA	NA	1 U [1 U]	1 U	2.8	1 U	1 U
t-Butyl Alcohol (TBA)	100	NA	NA	NA	NA	NA	NA	NA
1,1,2,2-Tetrachloroethane	1	NA	NA	0.33 J [0.32 J]	1 U	1 U	1 U	1 U
Tetrachloroethene	1	NA	NA	1 U [1 U]	1 U	1.3	1 U	0.96 J
Toluene	600	NA	NA	1 U [1 U]	1 U	0.52 J	1 U	1 U
1,1,1-Trichloroethane	30	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
1,1,2-Trichloroethane	3	NA	NA	1 U [1 U]	1 U	1 U	1 U	1 U
Trichloroethene	1	NA	NA	0.2 J [0.16 J]	0.29 J	6	0.63 J	69
Trichlorofluoromethane	2000	NA	NA	NA NA	NA	NA	NA	NA
Vinyl chloride	1	NA	NA	1 U [1 U]	1 U	1.9	1 U	1 U
Xylene (total)	1000	NA	NA	3 U [3 U]	3 U	3 U	3 U	3 U
Total VOCs	-	NA	NA	1.30 J [1.32 J]	0.91 J	209.47 J	1.26 J	108.65 J

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Attachment 2

2014 Baseline/PostInjection Event #1
Groundwater Analytical
Results

Volatile Organic Compounds (\	Sample ID Lab Sample ID Sample Date CRITERIA (7/22/2010)	EB JB57365-9	EB - FILTERED JB57365-9F	EB JB57365-19	EB - FILTERED	ТВ	EB	ТВ	EB	EB - FILTERED	ТВ	ТВ	ISCO-MW-1	ISCO-MW-1	ISCO-MW-2	ISCO-MW-2	ISCO-MW-3
Volatile Organic Compounds (\	Sample Date GROUNDWATER QUALITY  CRITERIA (7/22/2010)		JB57365-9F	IDE73CE 10									1300-111111-1	13CO-IVIVV-1			I3CO-IVIVV-3
Volatile Organic Compounds (\	Sample Date CRITERIA (7/22/2010)			102/302-13	JB57365-19F	JB57365-21	JB57510-11	JB57510-12	JB59106-2	JB59106-2F	JB59106-4	JB57131-4	JB57365-1	JB57365-1F	JB57365-7	JB57365-7F	JB57365-10
Volatile Organic Compounds (\		1/9/2014	1/9/2014	1/10/2014	1/10/2014	1/10/2014	1/13/2014	1/13/2014	2/3/2014	2/3/2014	2/3/2014	1/6/2014	1/9/2014	1/9/2014	1/10/2014	1/10/2014	1/9/2014
Volatile Organic Compounds (\	Matrix ug/L	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	GW	GW - FILTERED	GW	GW - FILTERED	GW
voiatile Organic Compounds (	Unit VOCs)	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
0	6000	ND (3.3)	<u>-</u>	ND (3.3)	_	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)		ND (3.3)	ND (3.3)	ND (3.3)		102		ND (3.3)
Acetone Benzene	1	ND (0.28)	-	ND (0.28)	-	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	0.83 J	-	ND (0.28)	-	ND (3.3) ND (0.28)
Bromochloromethane		ND (0.42)	-	ND (0.42)	-	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	-	ND (0.42)	ND (0.42)	ND (0.42)	-	ND (0.42)	-	ND (0.42)
Bromodichloromethane	1	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Bromoform	4	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Bromomethane	10	ND (0.56)	-	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)	-	ND (0.56)	-	ND (0.56)
2-Butanone (MEK)	300	ND (3.2)	-	ND (3.2)	-	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	-	ND (3.2)	ND (3.2)	ND (3.2)	-	173	-	ND (3.2)
Carbon disulfide	700	ND (0.18)	-	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	=	0.79	J -	ND (0.18)
Carbon tetrachloride	1	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	-	ND (0.23)	-	ND (0.23)
Chlorobenzene	50	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.35)	-	ND (0.35)
Chloroethane	70	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	-	ND (0.39) 1.7	-	ND (0.39) ND (0.25)
Chloroform Chloromethane	70	ND (0.36)	-	ND (0.36)	-	ND (0.25)	ND (0.25)	ND (0.36)	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	-	ND (0.36)	-	ND (0.36)
Cyclohexane		ND (0.36)	-	ND (0.18)	_	ND (0.38)	ND (0.38)	ND (0.18)	ND (0.38)	-	ND (0.38)	ND (0.38)	0.31 J	_	ND (0.38)	-	ND (0.18)
1,2-Dibromo-3-chloropropane	0.02	ND (1.3)	-	ND (1.3)	-	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	-	ND (1.3)	ND (1.3)	ND (1.3)	-	ND (1.3)	-	ND (1.3)
Dibromochloromethane	1	ND (0.19)	-	ND (0.19)	-	ND (0.19)	ND (0.19)	ND (0.19)	ND (0.19)	-	ND (0.19)	ND (0.19)	ND (0.19)	-	ND (0.19)	-	ND (0.19)
1,2-Dibromoethane	0.03	ND (0.16)	-	ND (0.16)	<u>-</u>	ND (0.16)	ND (0.16)	ND (0.16)	ND (0.16)	<u>-</u>	ND (0.16)	ND (0.16)	ND (0.16)	-	ND (0.16)	-	ND (0.16)
1,2-Dichlorobenzene	600	ND (0.20)	-	ND (0.20)	-	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	-	ND (0.20)	ND (0.20)	ND (0.20)	-	ND (0.20)	-	ND (0.20)
1,3-Dichlorobenzene	600	ND (0.31)	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	-	ND (0.31)	-	ND (0.31)
1,4-Dichlorobenzene	75	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Dichlorodifluoromethane	1000	ND (0.63)	-	ND (0.63)	-	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	-	ND (0.63)	ND (0.63)	ND (0.63)	-	ND (0.63)	-	ND (0.63)
1,1-Dichloroethane	50	ND (0.26)	-	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26)	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	ND (0.26)	0.46 J 21.4	-	ND (0.26)	-	ND (0.26)
1,2-Dichloroethane 1,1-Dichloroethene	1	ND (0.22) ND (0.34)	1 -	ND (0.34)	-	ND (0.22)	ND (0.22)	ND (0.22) ND (0.34)	ND (0.22)	-	ND (0.22)	ND (0.22) ND (0.34)	ND (0.34)	-	1270 ND (0.34)	-	0.48 J ND (0.34)
cis-1,2-Dichloroethene	70	ND (0.24)	<u> </u>	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.24)	ND (0.24)	3.2	-	1.8	<u> </u>	47.1
trans-1,2-Dichloroethene	100	ND (0.38)	-	ND (0.38)	-	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	-	ND (0.38)	ND (0.38)	ND (0.38)	-	3	-	ND (0.38)
1,2-Dichloropropane	1	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)
cis-1,3-Dichloropropene	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	ND (0.15)	ND (0.15)	ND (0.15)	-	ND (0.15)	ND (0.15)	ND (0.15)	-	ND (0.15)	-	ND (0.15)
trans-1,3-Dichloropropene	=	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)
1,4-Dioxane	-	ND (73)	-	ND (73)	-	ND (73)	ND (73)	ND (73)	ND (73)	-	ND (73)	ND (73)	ND (73)	-	ND (73)	-	ND (73)
Ethylbenzene	700	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Freon 113	=	ND (0.77)	-	ND (0.77)	-	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	-	ND (0.77)	ND (0.77)	ND (0.77)	-	ND (0.77)	-	ND (0.77)
2-Hexanone		ND (1.7)	-	ND (1.7)	-	ND (1.7)	ND (1.7)	ND (1.7)	ND (1.7)	-	ND (1.7)	ND (1.7)	ND (1.7)	-	ND (1.7)	-	ND (1.7)
Isopropylbenzene Methyl Acetate	700 7000	ND (0.22) ND (1.5)	-	ND (0.22) ND (1.5)	-	ND (0.22) ND (1.5)	ND (0.22) ND (1.5)	ND (0.22) ND (1.5)	ND (0.22) ND (1.5)	-	ND (0.22) ND (1.5)	ND (0.22) ND (1.5)	2.4 ND (1.5)	-	ND (0.22) ND (1.5)	-	ND (0.22) ND (1.5)
Methylcyclohexane	-	ND (0.15)	-	ND (0.15)	_	ND (0.15)	ND (0.15)	ND (0.15)	ND (0.15)	_	ND (0.15)	ND (0.15)	ND (0.15)	-	ND (0.15)	-	ND (0.15)
Methyl Tert Butyl Ether	70	ND (0.29)	-	ND (0.29)	-	ND (0.29)	ND (0.29)	ND (0.29)	ND (0.29)	-	ND (0.29)	ND (0.29)	ND (0.29)	-	ND (0.29)	-	ND (0.29)
4-Methyl-2-pentanone(MIBK)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	ND (1.5)	ND (1.5)	ND (1.5)	-	ND (1.5)	ND (1.5)	ND (1.5)	-	ND (1.5)	-	ND (1.5)
Methylene chloride	3	ND (0.86)	-	ND (0.86)	-	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	-	ND (0.86)	ND (0.86)	ND (0.86)	-	4.5	-	ND (0.86)
Styrene 1,1,2,2-Tetrachloroethane	100	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	-	ND (0.30) 22.5	-	ND (0.30) ND (0.20)
Tetrachloroethene	1	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	-	5	-	1.4
Toluene	600	ND (0.44)	-	ND (0.44)	-	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	-	ND (0.44)	ND (0.44)	ND (0.44)	-	1.7	-	ND (0.44)
1,2,3-Trichlorobenzene	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.24)	-	ND (0.24)
1,2,4-Trichlorobenzene 1,1,1-Trichloroethane	9 30	ND (0.22) ND (0.25)	-	ND (0.22) ND (0.25)	-	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	-	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) 0.46 J	-	ND (0.22) ND (0.25)	+	ND (0.22) ND (0.25)
1,1,2-Trichloroethane	3	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	0.49	J -	ND (0.23)
Trichloroethene	1	ND (0.50)	-	ND (0.50)		ND (0.50)	ND (0.50)	ND (0.50)	ND (0.50)	-	ND (0.50)	ND (0.50)	23.1	<u>-</u>	1.6	-	54.9
Trichlorofluoromethane	2000	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Vinyl chloride m,p-Xylene	1	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	-	ND (0.41) 0.43	-	ND (0.41) ND (0.40)
o-Xylene	-	ND (0.40)	-	ND (0.40)	-	ND (0.40) ND (0.19)	ND (0.40) ND (0.19)	ND (0.40)	ND (0.40)	-	ND (0.40)	ND (0.40)	ND (0.40) ND (0.19)	-	0.43	<u>,                                     </u>	ND (0.40) ND (0.19)
Xylene (total)	1000	ND (0.19)	-	ND (0.19)		ND (0.19)	ND (0.19)	ND (0.19)	ND (0.19)	<u>-</u>	ND (0.19)	ND (0.19)	ND (0.19)	<u>-</u>	0.67	J -	ND (0.19)
Total VOCs	-	0		0		0	0	0	0		0	0	52.16		1588.75		103.88
CC/MC V-I-Ail-TIC																	
GC/MS Volatile TIC Total TIC, Volatile	,	0 1		1 0 1 1		1 0 1	1 0 1	1 0 1	1 0 1	1	1 0 1	1 0 1	1 0 1	1 - 1	1 72		
Total TIC, Volatile Total Alkanes		0	-	0	-	0	0	0	0	-	0	0	0	-	7.2	J -	0
		-	ı L				<u>, , , , , , , , , , , , , , , , , , , </u>	, , , ,		1	,	<u> </u>	<u>, - 1</u>	1	<u>, , , , , , , , , , , , , , , , , , , </u>	1	
Metals Analysis																	الترسير
Chromium	70	<10	<10	<10	<10	<u> </u>	- 1	- 1	<10	<10	- 1	- 1	<10	<10	208		<10
Iron	300	- <10000	<100	- <10000	<100	-	-	-	- <10000	<100	-	-	18300	5670	22300	22200	18000
Sodium	50000	<10000	-	<10000	-	-	-	-	<10000	-	-	-	18300	-	22300	=	18000
General Chemistry																	
A STATE OF THE PARTY OF THE PAR	500000	<10000	- 1	<10000	- 1	- 1	] - 1	- 1	<10000	- 1	- 1		206000	-	433000	- 1	580000
Solids, Total Dissolved		<10000		<10000			1		<10000		-t		101000	-			241000

	Sample ID	ISCO-MW-3	ISCO-MW-3-DUP	ISCO-MW-3-DUP	ISCO-MW-4	ISCO-MW-4	ISCO-MW-5	ISCO-MW-5	ISCO-MW-6	ISCO-MW-6	ISCO-MW-7	ISCO-MW-7	ISCO-MW-8	ISCO-MW-8	ISCO-MW-9	ISCO-MW-9	IW1-BT-2
	Lab Sample ID	JB57365-10F	JB57365-11	JB57365-11F	JB59106-1	JB59106-1F	JB57365-3	JB57365-3F	JB57365-4	JB57365-4F	JB57365-18	JB57365-18F	JB57365-5	JB57365-5F	JB57365-20	JB57365-20F	JB57365-8
	Sample Date GROUNDWATER QUALITY	1/9/2014	1/9/2014	1/9/2014	2/3/2014	2/3/2014	1/9/2014	1/9/2014	1/9/2014	1/9/2014	1/10/2014	1/10/2014	1/9/2014	1/9/2014	1/10/2014	1/10/2014	1/9/2014
	Matrix (7/22/2010)	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW - FILTERED	GW
	ug/L Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (	(VOCs)																
Acetone	6000	-	ND (3.3)	-	ND (3.3)	-	ND (3.3)	-	ND (3.3)	-	ND (3.3)	-	30	-	ND (3.3)	-	ND (3.3)
Benzene	1	-	ND (0.28)	-	ND (0.28)	-	0.69 J	-	ND (0.28)	-	ND (0.28)	- +	ND (0.28)	-	ND (0.28)	-	ND (0.28)
Bromochloromethane Bromodichloromethane	- 1	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)	-	ND (0.42) ND (0.21)
Bromoform	4	-	ND (0.21)	-	ND (0.21) ND (0.30)	-	ND (0.30)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Bromomethane	10	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)
2-Butanone (MEK)	300	-	ND (3.2)	-	ND (3.2)	-	ND (3.2)	-	ND (3.2)	-	ND (3.2)	-	ND (3.2)	-	ND (3.2)	-	ND (3.2)
Carbon disulfide	700	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)
Carbon tetrachloride	1	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	0.8	J -	0.65	J -	ND (0.23)	-	ND (0.23)
Chlorobenzene	50	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)
Chloroethane	-	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)		ND (0.39)	-	ND (0.39)
Chloroform	70	-	ND (0.25)	-	0.36	J -	ND (0.25)	-	ND (0.25)	-	1.4	-	1.7	-	ND (0.25)	-	1.3
Chloromethane Cyclohexane	-	<del>-</del>	ND (0.36) ND (0.18)	+ -	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)
1,2-Dibromo-3-chloropropane	0.02	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	<del>                                     </del>	ND (0.18)	-	ND (0.18)	-	ND (0.18)
Dibromochloromethane	1	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)		ND (0.19)	-	ND (0.19)	-	ND (0.19)
1,2-Dibromoethane	0.03	-	ND (0.16)		ND (0.16)	- 1	ND (0.16)	-	ND (0.16)	-	ND (0.16)	- 1	ND (0.16)		ND (0.16)	-	ND (0.16)
1,2-Dichlorobenzene	600	-	ND (0.20)	<u> </u>	ND (0.20)	<u>-</u>	ND (0.20)	<u>-                                      </u>	ND (0.20)	=	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)
1,3-Dichlorobenzene	600	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)
1,4-Dichlorobenzene	75	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	=	ND (0.30)	÷	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Dichlorodifluoromethane	1000	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)
1,1-Dichloroethane 1,2-Dichloroethane	50 2	-	ND (0.26) 0.5	+	ND (0.26) 0.98	.1 -	ND (0.26) 46.7	-	ND (0.26) <b>0.56</b> J	-	ND (0.26) 1.1	-	ND (0.26) 36.8	-	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)
1,1-Dichloroethane	1	-	ND (0.34)		ND (0.34)	J -	ND (0.34)	-	ND (0.34)	-	ND (0.34)	1 -	ND (0.34)	-	ND (0.22)	-	ND (0.22)
cis-1,2-Dichloroethene	70	-	56.8	-	ND (0.24)	-	9.2	-	1.4	-	ND (0.24)	-	0.58	J -	2.6	-	ND (0.24)
trans-1,2-Dichloroethene	100	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)
1,2-Dichloropropane	1	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
cis-1,3-Dichloropropene	-	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)
trans-1,3-Dichloropropene	-	=	ND (0.21)	-	ND (0.21)	-	ND (0.21)	=	ND (0.21)	=	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
1,4-Dioxane	-	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)
Ethylbenzene	700	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Freon 113 2-Hexanone		-	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)	<del>                                     </del>	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)
Isopropylbenzene	700	-	ND (0.22)	-	ND (0.22)	-	0.61 J	_	ND (0.22)	-	ND (0.22)	-	ND (0.22)	<del>                                     </del>	ND (0.22)	-	ND (0.22)
Methyl Acetate	7000	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)
Methylcyclohexane	-	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)
Methyl Tert Butyl Ether	70	-	ND (0.29)	-	ND (0.29)	-	ND (0.29)	-	ND (0.29)	-	ND (0.29)	-	1.1	-	ND (0.29)	-	ND (0.29)
4-Methyl-2-pentanone(MIBK) Methylene chloride	- 2	-	ND (1.5) ND (0.86)	-	ND (1.5) 2.6	-	ND (1.5) 8.4	-	ND (1.5) ND (0.86)	-	ND (1.5) ND (0.86)	-	ND (1.5) 2.3	-	ND (1.5) ND (0.86)	-	ND (1.5) ND (0.86)
Styrene	100	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)
1,1,2,2-Tetrachloroethane	1	=	ND (0.20)	-	ND (0.20)	=	0.29 J	=	ND (0.20)	=	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)
Tetrachloroethene	1	-	1.6	-	ND (0.25)	-	0.79 J	-	ND (0.25)	-		J -	1.2	-		J -	ND (0.25)
Toluene 1,2,3-Trichlorobenzene	600	-	ND (0.44) ND (0.24)	+ -	ND (0.44) ND (0.24)	+	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	+	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)
1,2,4-Trichlorobenzene	9	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	+ - +	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,1,1-Trichloroethane	30	-	ND (0.25)	<u> </u>	ND (0.25)	<u> </u>	ND (0.25)		ND (0.25)	-	ND (0.25)		ND (0.25)	-	ND (0.25)	-	ND (0.25)
1,1,2-Trichloroethane	3	-	ND (0.21)		ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Trichloroethene Trichlorofluoromethane	1 2000	-	64 ND (0.33)	+ -	0.8 ND (0.33)	J -	17.1 ND (0.33)	-	2 ND (0.33)	-	ND (0.33)	+	5.8 ND (0.33)	-	44 ND (0.33)	-	1.6 ND (0.33)
Vinyl chloride	1	-	ND (0.41)		ND (0.41)	- 1	ND (0.41)	-	ND (0.41)	-	ND (0.41)	- +	ND (0.41)	-	ND (0.41)	-	ND (0.41)
m,p-Xylene	-	-	ND (0.40)	- 1	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)
o-Xylene	-	-	ND (0.19)		ND (0.19)		ND (0.19)	-	ND (0.19)	-	ND (0.19)	1	ND (0.19)	-	ND (0.19)	<del>                                     </del>	ND (0.19)
Xylene (total) Total VOCs	1000	-	ND (0.19) 122.9	-	ND (0.19) 4.74	+ -	ND (0.19) 83.78	-	ND (0.19) 3.96	-	ND (0.19) <b>6.21</b>	+	ND (0.19) 80.13	<del>                                     </del>	ND (0.19) 47.2	+ - +	ND (0.19) 2.9
	<u> </u>	<u> </u>	.22.0	1			00.70	1	5.50	I	V.E1		55.10				
GC/MS Volatile TIC																	
Total TIC, Volatile	-	-	0	-			0		0			- 1		-			0
Total Alkanes	-	-	0	-	0	-	0	-	0	-	0	-	0	-	0	-	0
Metals Analysis																	
Chromium	70	<10	<10	<10	16.4	<10	<10	<10	<10	<10	22.7	12.4	677	418	162	<10	<10
Iron	300		-	42600	-	<100	-		-	165	-	145	-	2920	-	1960	-
Sodium	50000	-	18800	-	15600	-	29000	-	54200	-	14300	- 1	599000	-	<10000	-	42000
Canada Chamista																	
General Chemistry Solids, Total Dissolved	500000		530000	- 1	46000		307000	1	297000		92000		1300000		203000	- 1	247000
Sulfate	250000	-	214000	-		-	154000	-	153000	-		-	682000	-	106000	-	F0000
		1					1	ı		1		1			1	1	1

Sample II	ID	IW1-BT-2	IW1-DR-1	IW1-DR-1	IW-4S	PZ-1S	PZ-1S	MW-5I	MW-5I	MW-6S	MW-9I	MW-9I-DUP	MW-10S	MW-10S	MW-11I	MW-11I	MW-14S
Lab Sample II		JB57365-8F	JB57365-12	JB57365-12F	JB57365-6	JB57365-2	JB57365-2F	JB57365-17	JB57365-17F	JB57510-2	JB57510-3	JB57510-4	JB57365-13	JB57365-13F	JB57365-16	JB57365-16F	JB57365-15
Sample Date	te CRITERIA (7/22/2010)	1/9/2014	1/10/2014	1/10/2014	1/9/2014	1/9/2014	1/9/2014	1/10/2014	1/10/2014	1/13/2014	1/13/2014	1/13/2014	1/10/2014	1/10/2014	1/10/2014	1/10/2014	1/10/2014
Matri	ix ug/L	GW - FILTERED	GW	GW - FILTERED	GW	GW	GW - FILTERED	GW	GW - FILTERED	GW	GW	GW	GW	GW	GW	GW	GW
Uni Volatile Organic Compounds (VOCs)	nit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
	5000	1	ND (0.0)	1	ND (0.0)	ND (3.3)	1	ND (0.0)		ND (0.0)	ND (0.0)	ND (0.0)	ND (0.0)	1	ND (0.0)	1	ND (2.2)
Acetone Benzene	6000	-	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)
Bromochloromethane	-	-	ND (0.42)	-	ND (0.42)	ND (0.42)	1 -	ND (0.42)	-	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	-	ND (0.42)	1 -	ND (0.42)
Bromodichloromethane	1	-	ND (0.21)	-	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Bromoform	4	-	ND (0.30)	-	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Bromomethane	10	-	ND (0.56)	-	ND (0.56)	ND (0.56)	-	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	-	ND (0.56)	-	ND (0.56)
2-Butanone (MEK)	300	-	ND (3.2)	-	ND (3.2)	ND (3.2)	=	ND (3.2)	=	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	=	ND (3.2)	=	ND (3.2)
Carbon disulfide	700	-	ND (0.18)	-	ND (0.18)	ND (0.18)	-	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	-	ND (0.18)	-	ND (0.18)
Carbon tetrachloride	1	-	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	-	ND (0.23)	-	ND (0.23)
Chlorobenzene	50	-	ND (0.35)	-	ND (0.35)	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.35)	-	ND (0.35)
Chloroethane	-	-	ND (0.39)	-	ND (0.39)	ND (0.39)	-	ND (0.39)	-	ND (0.39)	ND (0.39)	ND (0.39)	ND (0.39)	-	ND (0.39)	-	ND (0.39)
Chloroform	70	-	ND (0.25)	-	ND (0.25)	0.63	J -	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	0.62 J	-	0.27	J -	ND (0.25)
Chloromethane Cyclohexane	-	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	ND (0.36) ND (0.18)	ND (0.36) ND (0.18)	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)	-	ND (0.36) ND (0.18)
1,2-Dibromo-3-chloropropane	0.02	-	ND (1.3)	-	ND (0.18)	ND (0.18)	-	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	-	ND (1.3)	-	ND (0.18)
Dibromochloromethane	1	-	ND (0.19)	-	ND (0.19)	ND (0.19)	- 1	ND (0.19)	-	ND (0.19)	ND (0.19)	ND (0.19)	ND (0.19)	-	ND (0.19)	-	ND (0.19)
1,2-Dibromoethane	0.03	-	ND (0.16)	-	ND (0.16)	ND (0.16)	-	ND (0.16)	-	ND (0.16)	ND (0.16)	ND (0.16)	ND (0.16)	-	ND (0.16)	-	ND (0.16)
1,2-Dichlorobenzene	600	-	ND (0.20)	-	ND (0.20)	ND (0.20)	- 1	ND (0.20)	-	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	-	ND (0.20)	<u> </u>	ND (0.20)
1,3-Dichlorobenzene	600	-	ND (0.31)	-	ND (0.31)	ND (0.31)	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	-	ND (0.31)	-	ND (0.31)
1,4-Dichlorobenzene	75	-	ND (0.30)	-	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Dichlorodifluoromethane	1000	-	ND (0.63)	-	ND (0.63)	ND (0.63)	-	ND (0.63)	-	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	-	ND (0.63)	-	ND (0.63)
1,1-Dichloroethane	50	-	ND (0.26)	-	ND (0.26)	ND (0.26)		ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)
1,2-Dichloroethane	2	-	ND (0.22)	-	ND (0.22)	0.56	J -	1.9	-	ND (0.22)	ND (0.22)	ND (0.22)	1.2	-	ND (0.22)	-	ND (0.22)
1,1-Dichloroethene cis-1,2-Dichloroethene	70	-	ND (0.34) ND (0.24)	-	ND (0.34) 0.8	ND (0.34)	-	ND (0.34) ND (0.24)	-	ND (0.34) ND (0.24)	ND (0.34) ND (0.24)	ND (0.34) ND (0.24)	ND (0.34) 5.9	-	ND (0.34) ND (0.24)	-	ND (0.34) ND (0.24)
trans-1,2-Dichloroethene	100	-	ND (0.24)	-	ND (0.38)	ND (0.38)	-	ND (0.38)	-	ND (0.24)	ND (0.24)	ND (0.38)	ND (0.38)	-	ND (0.24)	-	ND (0.38)
1,2-Dichloropropane	100	-	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.28)	<del>-</del>	ND (0.28)
cis-1,3-Dichloropropene	-	-	ND (0.15)	-	ND (0.15)	ND (0.15)	-	ND (0.15)	-	ND (0.15)	ND (0.15)	ND (0.15)	ND (0.15)	-	ND (0.15)	-	ND (0.15)
trans-1,3-Dichloropropene	-	-	ND (0.21)	-	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)
1,4-Dioxane	-	-	ND (73)	-	ND (73)	ND (73)	-	ND (73)	-	ND (73)	ND (73)	ND (73)	ND (73)	-	ND (73)	-	ND (73)
Ethylbenzene	700	-	ND (0.21)	-	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Freon 113	-	-	ND (0.77)	-	ND (0.77)	ND (0.77)	-	ND (0.77)	-	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	-	ND (0.77)	-	ND (0.77)
2-Hexanone	-	-	ND (1.7)	-	ND (1.7)	ND (1.7)	-	ND (1.7)	-	ND (1.7)	ND (1.7)	ND (1.7)	ND (1.7)	-	ND (1.7)	-	ND (1.7)
Isopropylbenzene	700	-	ND (0.22)	-	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)
Methyl Acetate Methylcyclohexane	7000	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	ND (1.5) ND (0.15)	ND (1.5) ND (0.15)	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)
Methyl Tert Butyl Ether	70	-	ND (0.29)	-	ND (0.19)	ND (0.19)	-	ND (0.29)	-	ND (0.29)	ND (0.29)	ND (0.29)	ND (0.29)	-	ND (0.19)	-	ND (0.29)
4-Methyl-2-pentanone(MIBK)	-	-	ND (1.5)	-	ND (1.5)	ND (1.5)	-	ND (1.5)	-	ND (1.5)	ND (1.5)	ND (1.5)	ND (1.5)	-	ND (1.5)	-	ND (1.5)
Methylene chloride	3	-	ND (0.86)	-	ND (0.86)	ND (0.86)	-	ND (0.86)	-	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	-	ND (0.86)	-	ND (0.86)
Styrene 1,1,2,2-Tetrachloroethane	100	-	ND (0.30) ND (0.20)	+	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)
Tetrachloroethene	1	-	0.72	J -	ND (0.25)	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	ND (0.25)	-	ND (0.25)	-	ND (0.25)
Toluene	600	-	ND (0.44)	-	ND (0.44)	ND (0.44)		ND (0.44)	-	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	-	ND (0.44)	-	ND (0.44)
1,2,3-Trichlorobenzene	-	-	ND (0.24)	-	ND (0.24)	ND (0.24)	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.24)	-	ND (0.24)
1,2,4-Trichlorobenzene	9	-	ND (0.22)	-	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	ND (0.22)	-	ND (0.22)		ND (0.22)
1,1,1-Trichloroethane 1,1,2-Trichloroethane	30	-	ND (0.25) ND (0.21)	-	ND (0.25) ND (0.21)	ND (0.25) ND (0.21)	-	ND (0.25) ND (0.21)	-	0.42 ND (0.21)	J ND (0.25) ND (0.21)	ND (0.25) ND (0.21)	ND (0.25) ND (0.21)	-	1.6 ND (0.21)	-	ND (0.25) ND (0.21)
Trichloroethene	1	-	1.1	-	2.4	2	-	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)	5.5	-	ND (0.50)	-	ND (0.50)
Trichlorofluoromethane	2000	-	ND (0.33)	-	ND (0.33)	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Vinyl chloride	1	-	ND (0.41)	-	ND (0.41)	ND (0.41)	-	ND (0.41)	-	ND (0.41)	ND (0.41)	ND (0.41)	ND (0.41)	-	ND (0.41)	-	ND (0.41)
m,p-Xylene o-Xylene	-	-	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	ND (0.40) ND (0.19)	+	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	ND (0.40) ND (0.19)	ND (0.40) ND (0.19)	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)
Xylene (total)	1000	-	ND (0.19)	<del>-</del>	ND (0.19)	ND (0.19)	-	ND (0.19)	-	ND (0.19)	ND (0.19)		ND (0.19)		ND (0.19)	-	ND (0.19)
Total VOCs			1.82		3.2	4.59		1.9		0.42	0	0	13.22		1.87		
GC/MS Volatile TIC							,										
Total TIC, Volatile Total Alkanes	-	-	0	-	0 0		-	0		0	0	0	0	-	0	-	0
TOTAL AIRAITES	<u> </u>	1 -	U		U U	U	<u> </u>	U	-	U	U	U	U	-	ı	<u> </u>	
Metals Analysis																	
Chromium	70	<10	<10	<10	-	<10	<10	<10		- 1	- 1	-	<10		<10	<10	<10
Iron	300	235	-	27500	-		<100	-	134	-	-	-	-	<100	-	<100	
Sodium	50000	-	12200	-	-	16500	-	<10000	-	-	-	-	25500	-	<10000	-	229000
General Chemistry																	
Solids, Total Dissolved	500000	- 1	218000	-	- 1	384000	- 1	205000	- 1	- 1	- 1		214000	-	118000	- 1	1090000
Sulfate	250000	-	101000	-	-	177000	-	68000	-	- 1	-	-	84800	-	39900	-	
	•					·		•			•			•		_ ·	<del></del>

Sample ID	)	MW-14S	MW-14SD	MW-14SD	MW-15D	MW-19S	MW-23S	MW-23I	MW-23D	MW-24	MW-28	WCC-1S	WCC-1M	WCC-3M	EW-3	EB-01_05122014	EB-01_05122014
Lab Sample ID		JB57365-15F	JB57365-14	JB57365-14F	JB57510-9	JB57510-7	JB57131-2	JB57131-3	JB57131-1	JB57510-1	JB57510-8	JB57510-6	JB57510-5	JB57510-10	JB59106-3	JB66824-5	JB66824-5F
Sample Date	CRITERIA (7/22/2010)	1/10/2014	1/10/2014	1/10/2014	1/13/2014	1/13/2014	1/6/2014	1/6/2014	1/6/2014	1/13/2014	1/13/2014	1/13/2014	1/13/2014	1/13/2014	2/3/2014	5/12/2014	5/12/2014
Matrix Unit	ug/L	GW	GW	GW ug/L	GW	GW	GW	GW	GW ug/L	GW	GW	GW	GW	GW ug/L	GW	WATER	WATER-FILTERED
Volatile Organic Compounds (VOCs)		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Acetone	6000	-	ND (3.3)	=	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	ND (3.3)	-
Benzene	1	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	ND (0.28)	=
Bromochloromethane	-	-	ND (0.42)	-	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	ND (0.42)	=
Bromodichloromethane Bromoform	1 4	-	ND (0.21) ND (0.30)	-	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	ND (0.21) ND (0.30)	=
Bromomethane	10	-	ND (0.56)	-	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.56)	ND (0.30) ND (0.56)	-
2-Butanone (MEK)	300	-	ND (3.2)	-	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	ND (3.2)	<u>-</u>
Carbon disulfide	700	-	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	=
Carbon tetrachloride	1	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	ND (0.23)	-
Chloroethane	50	-	ND (0.35) ND (0.39)	-	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35) ND (0.39)	ND (0.35)	-
Chloroform	70	-	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	1.6	ND (0.25)	ND (0.39)	ND (0.39)	ND (0.39) ND (0.25)	ND (0.39)	0.46	J 0.5 J	ND (0.39)	ND (0.39) ND (0.25)	-
Chloromethane	-	-	ND (0.36)	-	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	ND (0.36)	-
Cyclohexane	-	-	ND (0.18)	-	ND (0.18)	1.1 J	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	ND (0.18)	-
1,2-Dibromo-3-chloropropane	0.02	-	ND (1.3)	-	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	ND (1.3)	-
Dibromochloromethane 1,2-Dibromoethane	0.03	-	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	-
1,2-Dichlorobenzene	600	-	ND (0.10)	-	ND (0.10)	ND (0.10)	ND (0.10)	ND (0.20)	ND (0.10)	ND (0.10)	ND (0.10)	ND (0.10)	ND (0.20)	ND (0.10)	ND (0.20)	ND (0.16)	-
1,3-Dichlorobenzene	600	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	ND (0.31)	-
1,4-Dichlorobenzene	75	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	-
Dichlorodifluoromethane	1000	-	ND (0.63)	-	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	-
1,1-Dichloroethane 1,2-Dichloroethane	50 2	-	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	0.42 J ND (0.22)	ND (0.26) 15.6	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26) 35.1	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	-
1,1-Dichloroethene	1	-	ND (0.34)	-	2	ND (0.34)	ND (0.34)	ND (0.34)	ND (0.34)	ND (0.34)	ND (0.34)	ND (0.34)	ND (0.34)	0.43 J	ND (0.34)	ND (0.22)	-
cis-1,2-Dichloroethene	70	-	ND (0.24)	-	ND (0.24)	5.4	5.1	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	7.7	ND (0.24)	ND (0.24)	ND (0.24)	-
trans-1,2-Dichloroethene	100	-	ND (0.38)	-	ND (0.38)	1.8	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	ND (0.38)	-
1,2-Dichloropropane	1	-	ND (0.28) ND (0.15)	-	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28)	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28)	-
cis-1,3-Dichloropropene trans-1,3-Dichloropropene	-	-	ND (0.15) ND (0.21)	-	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	ND (0.15) ND (0.21)	-
1,4-Dioxane	-	-	ND (73)	-	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	ND (73)	-
Ethylbenzene	700	-	ND (0.21)	-	ND (0.21)	5.3	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-
Freon 113	-	-	ND (0.77)	-	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	ND (0.77)	-
2-Hexanone Isopropylbenzene	700	-	ND (1.7) ND (0.22)	-	ND (1.7) ND (0.22)	ND (1.7) 0.77 J	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	ND (1.7) 0.28 J	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	ND (1.7) ND (0.22)	-
Methyl Acetate	7000	-	ND (0.22)	-	ND (1.5)	ND (1.5)	ND (1.5)	ND (1.5)	ND (0.22)	ND (1.5)	ND (1.5)	ND (0.22)	ND (1.5)	ND (0.22)	ND (0.22)	ND (0.22) ND (1.5)	-
Methylcyclohexane	-	-	ND (0.15)	-	ND (0.15)	1.8 J	ND (0.15)	ND (0.15)	ND (0.15)	ND (0.15)	0.4 J	ND (0.15)	ND (0.15)	ND (0.15)	ND (0.15)	ND (0.15)	-
Methyl Tert Butyl Ether	70	-	ND (0.29) ND (1.5)	-	ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	0.63 ND (1.5)	J ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	1.3 ND (1.5)	ND (0.29) ND (1.5)	ND (0.29) ND (1.5)	-
4-Methyl-2-pentanone(MIBK) Methylene chloride	3	-	ND (0.86)	-	ND (0.86)	ND (0.86)	4.7	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	ND (0.86)	ND (1.5)	-
Styrene	100	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	ND (0.30)	-
1,1,2,2-Tetrachloroethane Tetrachloroethene	1 1	-	ND (0.20) ND (0.25)	-	ND (0.20) ND (0.25)	ND (0.20) ND (0.25)	0.36 0.27	J ND (0.20) J ND (0.25)	ND (0.20) ND (0.25)	ND (0.20) ND (0.25)	ND (0.20) 0.27 J	ND (0.20) ND (0.25)	0.29 0.27	J 0.23 J J ND (0.25)	0.48 ND (0.25)	J ND (0.20) ND (0.25)	-
Toluene	600	-	ND (0.44)		ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.44)	ND (0.23)	-
1,2,3-Trichlorobenzene	-	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	ND (0.24)	-
1,2,4-Trichlorobenzene 1,1,1-Trichloroethane	9 30	-	ND (0.22) ND (0.25)		ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	ND (0.22) ND (0.25)	-
1,1,2-Trichloroethane	3	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	ND (0.21)	-
Trichloroethene	1	-	ND (0.50) ND (0.33)	-	ND (0.50)	4 ND (0.33)	7.3 ND (0.33)	ND (0.50) ND (0.33)	ND (0.50)	ND (0.50) ND (0.33)	ND (0.50) ND (0.33)	ND (0.50) ND (0.33)	5.1 ND (0.33)	ND (0.50) ND (0.33)	ND (0.50) ND (0.33)	ND (0.50)	-
Trichlorofluoromethane Vinyl chloride	2000	-	ND (0.33) ND (0.41)	-	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	ND (0.33) ND (0.41)	-
m,p-Xylene	-	-	ND (0.40)	-	ND (0.40)	1.3	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	ND (0.40)	-
o-Xylene Xylene (total)	1000	-	ND (0.19) ND (0.19)	-	ND (0.19) ND (0.19)	1.3 2.5	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	ND (0.19) ND (0.19)	-
Total VOCs	-		0	· ·	2	23.09	(/	0	0.63	0	( 7	0	48.92	2.46	0.48	0 ND (0.19)	-
	•					•	•						•				· ·
GC/MS Volatile				,		175.0					150.0		77			1 22	1
Total TIC, Volatile Total Alkanes	-	-	0	-	0			0		0		0 0	7.7	J 0	0		J -
											· · · · · · · · · · · · · · · · · · ·	· · · · · · · · · · · · · · · · · · ·			· · · · · ·		
Metals Analysis																	
Chromium Iron	70 300	<10 <100	<10	<10 <100	-	-	-	-	-	-		-	-	-	-	<10	<10 <100
Sodium	50000		228000	-	-		-	-	-	-			-	-	-		-
General Chemistry Solids, Total Dissolved	500000	_	1440000		1 -						1 - 1	1 - 1				<10000	
Sulfate	250000	-	394000		+	-	-	-	-	-	-	<del>                                     </del>	<del>-</del>	-	-	<10000	-
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	Sample ID NJ CLASS IIA	EB-02_05132014	EB-02_05132014	EB-03_05142014	EB-03_05142014	TB-01_05132014	TB-02 _05142014	1RND1_ISCOMW-1	1RND1_ISCOMW-1	1RND1_ISCOMW-2	1RND1_ISCOMW-2	1RND1_ISCOMW-3	1RND1_ISCOMW-3	1RND1_ISCOMW-4	1RND1_ISCOMW-4	1RND1_ISCOMW-5
	Sample ID GROUNDWATER QUALITY ample Date	JB66824-9 5/13/2014	JB66824-9F 5/13/2014	JB66824-19 5/14/2014	JB66824-19F 5/14/2014	JB66824-10 5/13/2014	JB66824-20 5/14/2014	JB66824-14 5/14/2014	JB66824-14F 5/14/2014	JB66824-15 5/14/2014	JB66824-15F 5/14/2014	JB66824-4 5/12/2014	JB66824-4F 5/12/2014	JB66824-2 5/12/2014	JB66824-2F 5/12/2014	JB66824-16 5/14/2014
34	Matrix (7/22/2010)	WATER	WATER-FILTERED	WATER	WATER-FILTERED	WATER	WATER	GW	GW-FILTERED	3/14/2014 GW	GW-FILTERED	3/12/2014 GW	GW-FILTERED	3/12/2014 GW	GW-FILTERED	3/14/2014 GW
	ug/L Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs		ND (2.2)		ND (2.2)		ND (2.2)	ND (2.2)		1	455			1	ND (2.2)	1	40.4
Acetone Benzene	6000 1	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	ND (3.3) ND (0.28)	6.9 J ND (0.28)	-	177 0.47 J	-	94.4 ND (0.28)	-	ND (3.3) ND (0.28)	-	13.4 0.67
Bromochloromethane	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	ND (0.42)	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)
Bromodichloromethane	1	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Bromoform	4	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Bromomethane 2-Butanone (MEK)	10 300	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)	ND (0.56) ND (3.2)	ND (0.56) ND (3.2)	-	0.87 J 23.5	-	2 ND (3.2)	-	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)
Carbon disulfide	700	ND (0.18)	-	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	-	0.59 J	-	0.28	J -	ND (0.18)	-	ND (0.18)
Carbon tetrachloride	1	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	-	2	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)
Chlorobenzene	50	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)
Chloroethane Chloroform	70	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	ND (0.39) ND (0.25)	-	ND (0.39) <b>4.7</b>	-	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)	-	ND (0.39) ND (0.25)
Chloromethane	-	ND (0.36)	-	ND (0.36)	-	ND (0.36)	ND (0.36)	ND (0.36)	-	ND (0.36)	-	ND (0.36)	-	ND (0.36)	-	ND (0.36)
Cyclohexane	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	ND (0.18)	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)
1,2-Dibromo-3-chloropropane	0.02	ND (1.3)	-	ND (1.3)	-	ND (1.3)	ND (1.3)	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)
Dibromochloromethane 1,2-Dibromoethane	0.03	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)	-	ND (0.19) ND (0.16)
1,2-Dichlorobenzene	600	ND (0.20)	-	ND (0.20)		ND (0.20)	ND (0.20)	ND (0.20)	<u>-</u>	0.21 J	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)
1,3-Dichlorobenzene	600	ND (0.31)	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)
1,4-Dichlorobenzene	75	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)
Dichlorodifluoromethane 1,1-Dichloroethane	1000 50	ND (0.63) ND (0.26)	-	ND (0.63) ND (0.26)	-	ND (0.63) ND (0.26)	ND (0.63) ND (0.26)	ND (0.63) ND (0.26)	-	ND (0.63) ND (0.26)	-	ND (0.63) ND (0.26)	-	ND (0.63) ND (0.26)	-	ND (0.63) <b>0.45</b> J
1,2-Dichloroethane	2	ND (0.22)	-	ND (0.20)	-	ND (0.20)	ND (0.22)	5.9	-	21.3	-	0.38	J -	ND (0.22)	-	30.9
1,1-Dichloroethene	1	ND (0.34)	-	ND (0.34)	-	ND (0.34)	ND (0.34)	ND (0.34)	-	ND (0.34)	-	ND (0.34)	-	ND (0.34)	-	ND (0.34)
cis-1,2-Dichloroethene	70	ND (0.24)	-	ND (0.24)	-	ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.24)	-	39	-	ND (0.24)	-	1.8
trans-1,2-Dichloroethene 1,2-Dichloropropane	100	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)	ND (0.38) ND (0.28)	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)
cis-1,3-Dichloropropene	-	ND (0.28)	-	ND (0.28)	-	ND (0.28) ND (0.15)	ND (0.28) ND (0.15)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
trans-1,3-Dichloropropene	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
1,4-Dioxane	-	ND (73)	-	ND (73)	-	ND (73)	ND (73)	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (73)
Ethylbenzene Freon 113	700	ND (0.21) ND (0.77)	-	ND (0.21) ND (0.77)	-	ND (0.21) ND (0.77)	ND (0.21) ND (0.77)	ND (0.21) ND (0.77)	-	ND (0.21) ND (0.77)	-	ND (0.21) ND (0.77)	-	ND (0.21) ND (0.77)	-	ND (0.21) ND (0.77)
2-Hexanone	-	ND (0.77)	-	ND (0.77) ND (1.7)	-	ND (0.77) ND (1.7)	ND (0.77)	ND (0.77)	-	ND (0.77)	-	ND (1.7)	-	ND (1.7)	-	ND (0.77)
Isopropylbenzene	700	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	0.74 J
Methyl Acetate	7000	ND (1.5)	-	ND (1.5)	-	ND (1.5)	ND (1.5)	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)
Methylcyclohexane Methyl Tert Butyl Ether	70	ND (0.15) ND (0.29)	-	ND (0.15) ND (0.29)	-	ND (0.15) ND (0.29)	ND (0.15) ND (0.29)	ND (0.15) ND (0.29)	-	ND (0.15) ND (0.29)	-	ND (0.15) ND (0.29)	-	ND (0.15) ND (0.29)	-	ND (0.15) ND (0.29)
4-Methyl-2-pentanone(MIBK)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	ND (1.5)	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)
Methylene chloride Styrene	3 100	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	ND (0.86) ND (0.30)	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)
1,1,2,2-Tetrachloroethane	1	ND (0.20)	-	ND (0.30) ND (0.20)	-	ND (0.30) ND (0.20)	ND (0.20)	ND (0.20)	-	50.1	-	ND (0.30)	-	ND (0.30)	-	ND (0.20)
Tetrachloroethene	1	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	-	16.6	-		J -	ND (0.25)	-	ND (0.25)
Toluene 1,2,3-Trichlorobenzene	600	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	ND (0.44) ND (0.24)	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)		ND (0.44) ND (0.24)
1,2,4-Trichlorobenzene	9	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	ND (0.22)	-	1.3 J	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,1,1-Trichloroethane 1,1,2-Trichloroethane	30	ND (0.25) ND (0.21)	-	ND (0.25) ND (0.21)	-	ND (0.25) ND (0.21)	ND (0.25) ND (0.21)	ND (0.25) ND (0.21)	-	ND (0.25) 0.72 J	-	ND (0.25) ND (0.21)	-	ND (0.25) ND (0.21)	-	0.54 J ND (0.21)
Trichloroethene	1	ND (0.21) ND (0.50)	-	ND (0.21)	<u>-</u>	ND (0.21)	ND (0.21) ND (0.50)	0.72 J	<u>-</u>	60.7	-	33.2	-	ND (0.21) ND (0.50)	-	22.2
Trichlorofluoromethane	2000	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Vinyl chloride m,p-Xylene	- 1	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)	-	ND (0.41) ND (0.40)		ND (0.41) ND (0.40)
o-Xylene	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	ND (0.19)	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	1.1
Xylene (total) Total VOCs	1000	ND (0.19) 0	-	ND (0.19) 0	-	ND (0.19) 0	ND (0.19) <b>0</b>	ND (0.19) 13.52	-	ND (0.19) <b>360.06</b>	-	ND (0.19) 170.01	-	ND (0.19) <b>0</b>	-	1.1 71.8
Total VOCS	-	U					· · ·	13.32	<u> </u>	300.00	<u> </u>	170.01		1 0	1	/1.0
GC/MS Volatile TIC					_											
Total TIC, Volatile Total Alkanes	-	0	-	0	-	0	0	0	-	0	-	0	-	0	-	0
TOTAL AIRCITES	<u> </u>	U	<u> </u>	<u> </u>		<u> </u>	· · ·	U	<u> </u>	U	<u> </u>	U				U
Metals Analysis																
Chromium	70 300	<10	<10 <100	<10		-	-	28.6	<10 <b>236</b>			11700	8990 <100	<10		<10
Iron Sodium	300 50000	<10000	<100	<10000	1200	-	-	97000	- 236		238	<100000	<100	13400	<100	41200
General Chemistry	500000	×10000	,	<10000	- 1		, , ,	136000	- 1	8080000	,	14200000		35000	,	241000
Solids, Total Dissolved Sulfate	500000 250000	<10000 <10000	-	<10000 <10000		-		136000 137000		8080000 2020000	-			35000 36000		241000 95100
						ı					1		I.		1	

Sample ID	NJ CLASS IIA	1RND1_ISCOMW-5 JB66824-16F	1RND1_DUP01 JB66824-18	1RND1_DUP01 JB66824-18F	1RND1_ISCOMW-6 JB66824-17	1RND1_ISCOMW-6 JB66824-17F	1RND1_ISCOMW-7 JB66824-3	1RND1_ISCOMW-7 JB66824-3F	1RND1_ISCOMW-8 JB66824-1	1RND1_ISCOMW-8 JB66824-1F	1RND1_ISCOMW-9 JB66824-13	1RND1_ISCOMW-9 JB66824-13F	1RND1_IW1-BT-2 JB66824-12	1RND1_IW1-BT-2 JB66824-12F	1RND1_MW-10S JB66824-8	1RND1_MW-10S JB66824-8F
Lab Sample ID Sample Date	GROUNDWATER QUALITY	5/14/2014	5/14/2014	5/14/2014	5/14/2014	5/14/2014	5/12/2014	5/12/2014	5/12/2014	5/12/2014	5/13/2014	5/13/2014	5/13/2014	5/13/2014	5/13/2014	5/13/2014
Matrix		GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED
Unit		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs)	6000	_	ND (3.3)	1	ND (3.3)		26.6		26.1	ı	36.7		17.5	1	ND (2.2)	
Acetone Benzene	6000 1	-	0.66	-	ND (3.3)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (3.3) ND (0.28)	-
Bromochloromethane	-	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-	ND (0.42)	-
Bromodichloromethane	1	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-
Bromoform	4	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-
Bromomethane 2-Butanone (MEK)	10 300	-	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)	-	ND (0.56) ND (3.2)	-	1.1 ND (3.2)	-	ND (0.56) ND (3.2)	-
Carbon disulfide	700	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	0.32 J	-	0.5	J -		-	ND (0.18)	-
Carbon tetrachloride	1	-	ND (0.23)	-	ND (0.23)	-	0.81 J	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-
Chlorobenzene	50	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	=,	ND (0.35)		ND (0.35)	=,
Chloroform	- 70	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39) 0.38	-	ND (0.39)	-
Chloroform Chloromethane	70	-	ND (0.25) ND (0.36)	-	ND (0.25) ND (0.36)	-	1.5 ND (0.36)	-	0.83 J ND (0.36)	-	ND (0.25) ND (0.36)	-			ND (0.25) ND (0.36)	-
Cyclohexane	-	-	ND (0.18)	<u>-</u>	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-
1,2-Dibromo-3-chloropropane	0.02	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-	ND (1.3)	-
Dibromochloromethane	1	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-
1,2-Dibromoethane 1,2-Dichlorobenzene	0.03 600	-	ND (0.16) ND (0.20)	-	ND (0.16) ND (0.20)	-	ND (0.16)	-	ND (0.16)	-	ND (0.16)	=	ND (0.16) ND (0.20)	-	ND (0.16) ND (0.20)	-
1,2-Dichlorobenzene 1,3-Dichlorobenzene	600	-	ND (0.20) ND (0.31)	<del>                                     </del>	ND (0.20) ND (0.31)	-	ND (0.20) ND (0.31)	-	ND (0.20) ND (0.31)	-	ND (0.20) ND (0.31)	-	ND (0.20) ND (0.31)		ND (0.20) ND (0.31)	-
1,4-Dichlorobenzene	75	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	=	ND (0.30)	=	ND (0.30)	=
Dichlorodifluoromethane	1000	÷	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	-	ND (0.63)	=	ND (0.63)	=	ND (0.63)	=
1,1-Dichloroethane	50	-	0.48	J -	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	=	ND (0.26)	=	ND (0.26)	=
1,2-Dichloroethane 1,1-Dichloroethene	2	-	<b>32.2</b> ND (0.34)	-	2.2 ND (0.34)	-	ND (0.22) ND (0.34)	-	1 ND (0.34)	-	ND (0.22) ND (0.34)	-	ND (0.22) ND (0.34)	-	ND (0.22) ND (0.34)	-
cis-1,2-Dichloroethene	70	-	1.8	-	2.5	-	ND (0.34) ND (0.24)	-	0.54 J	-	0.91	1 -	ND (0.34)	-	ND (0.34) ND (0.24)	-
trans-1,2-Dichloroethene	100	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	-	ND (0.38)	=	ND (0.38)	=
1,2-Dichloropropane	1	÷	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	=	ND (0.28)	=	ND (0.28)	=
cis-1,3-Dichloropropene	-	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-	ND (0.15)	-
trans-1,3-Dichloropropene 1,4-Dioxane	-	-	ND (0.21) ND (73)	-	ND (0.21) ND (73)	-	ND (0.21) ND (73)	-	ND (0.21) ND (73)	-	ND (0.21) ND (73)	-	ND (0.21) ND (73)	-	ND (0.21) ND (73)	-
Ethylbenzene	700	-	ND (73) ND (0.21)	-	ND (73) ND (0.21)	-	ND (73) ND (0.21)	-	ND (0.21)	-	ND (73) ND (0.21)	-	ND (73) ND (0.21)	-	ND (73) ND (0.21)	-
Freon 113	-	-	ND (0.77)	-	ND (0.77)	-	ND (0.77)	-	ND (0.77)	-	ND (0.77)	=	ND (0.77)	=	ND (0.77)	=
2-Hexanone	-	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-
Isopropylbenzene	700	-	0.74 .	J -	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-
Methyl Acetate  Methylcyclohexane	7000	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-	ND (1.5) ND (0.15)	-
Methyl Tert Butyl Ether	70	-	ND (0.13)	-	ND (0.13)	-	ND (0.13) ND (0.29)	-	ND (0.13) ND (0.29)	-	ND (0.13)	=	ND (0.13)	=	ND (0.13)	-
4-Methyl-2-pentanone(MIBK)	-	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-
Methylene chloride Styrene	3 100	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	=	ND (0.86) ND (0.30)	=	ND (0.86) ND (0.30)	-
1,1,2,2-Tetrachloroethane	1	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	=	ND (0.20)	=	ND (0.20)	=
Tetrachloroethene	1	-	ND (0.25)	-	0.36 J	-	0.83 J	-	0.63 J	-	0.55	J -	ND (0.25)	-	ND (0.25)	-
Toluene 1.2.3-Trichlorobenzene	600	-	ND (0.44) ND (0.24)	+ -	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-	ND (0.44) ND (0.24)	-
1,2,4-Trichlorobenzene	9	-	ND (0.22)	<u> -                                   </u>	ND (0.22)		ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	<u>-</u>	ND (0.22)	-
1,1,1-Trichloroethane	30	-	0.54	J -	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-
1,1,2-Trichloroethane Trichloroethene	3 1	-	ND (0.21) 21.9	-	ND (0.21) 6.1	-	ND (0.21)	-	ND (0.21) 2.4	-	ND (0.21) 21.1	=	ND (0.21) 3.1	-	ND (0.21) ND (0.50)	-
Trichlorofluoromethane	2000		ND (0.33)	<u>-</u>	ND (0.33)		ND (0.33)		ND (0.33)		ND (0.33)		ND (0.33)	<u>-</u>	ND (0.33)	
Vinyl chloride	1	-	ND (0.41)	-	ND (0.41)	-	ND (0.41)	-	ND (0.41)	-	ND (0.41)	-	ND (0.41)	-	ND (0.41)	-
m,p-Xylene o-Xylene	-	-	ND (0.40) 1.2	<del> </del>	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	-	ND (0.40) ND (0.19)	-
Xylene (total)	1000	-	1.2	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-
Total VOCs	-		59.52		11.16		32.74		31.82		59.76		24.52		0	
GC/MS Volatile TIC																
Total TIC, Volatile	- 1	- 1	0		0	- 1	0	- 1	0	- 1	0	- 1	0	- 1	0	- 1
Total Alkanes	-	-	0	-	0	-	0	-	0	-	0	-	0	-	0	-
Metals Analysis																
Chromium	70	<10	<10	<10	<10	<10	946	1060	5560	6320	538 a	220	863 a	338	<10	<10
Iron	300	101	-	<100	-	<100	-	17600	-	2360	-	6650	-	1220	-	<100
Sodium	50000	-	40600	-	48100	-	784000	-	2250000	-	578000 a	=	344000 a	-	<10000	=
General Chemistry																
Solids, Total Dissolved	500000	-	227000	- 1		-	4430000	- 1	10900000	-	2430000	-	1600000	-		-
Sulfate	250000	-	94100	-	134000	-	511000	-	2600000	-	742000	-	518000	-	22200	-

Sample ID		1RND1 MW-14SA	1RND1 MW-14SA	1RND1 MW-14SB	1RND1 MW-14SB	1RND1 PZ-1S	1RND1 PZ-1S	1RND2 FB 06302014	1RND2 FB 06302014	1RND2 FB 070120	14 1RND2 FB 07012014	ТВ	ТВ	1RND2 ISCO-MW1	1RND2 ISCO-MW1	1RND2 ISCO-MW2
Lab Sample ID	NICIASSIIA	JB66824-6	JB66824-6F	JB66824-7	JB66824-7F	JB66824-11	JB66824-11F	JB70619-6	JB70619-6F	JB70619-16	JB70619-16F	JB70619-9	JB70619-17	JB70619-3	JB70619-3F	JB70619-19
Sample Date	CRITERIA (7/22/2010)	5/13/2014	5/13/2014	5/13/2014	5/13/2014	5/13/2014	5/13/2014	6/30/2014	6/30/2014	7/1/2014	7/1/2014	7/1/2014	7/1/2014	6/30/2014	6/30/2014	7/1/2014
Matrix	ug/L	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	WATER	WATER-FILTERED	WATER	WATER-FILTERED	WATER	WATER	GW	GW-FILTERED	GW
Unit		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs)	5000	ND (2.2)	1 1	ND (2.2)	ı		1	l wa (2.5)	1	ND (2.5)		ND (2.5)	ND (2.6)	ND (2.5)	1	1 204
Acetone Benzene	6000	ND (3.3) ND (0.28)	-	ND (3.3) ND (0.28)	-	18.1 ND (0.28)	-	ND (2.6) ND (0.21)	-	ND (2.6) ND (0.21)	-	ND (2.6) ND (0.21)	ND (2.6) ND (0.21)	ND (2.6) ND (0.21)	-	291 0.64 J
Bromochloromethane	-	ND (0.42)		ND (0.42)	_	ND (0.42)	-	ND (0.49)	-	ND (0.21)	_	ND (0.49)	ND (0.49)	ND (0.49)	-	ND (1.2)
Bromodichloromethane	1	ND (0.21)	- 1	ND (0.21)	-	ND (0.21)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	-	ND (0.70)
Bromoform	4	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	ND (0.31)	ND (0.31)	-	ND (0.78)
Bromomethane	10	ND (0.56)	E	ND (0.56)	=	ND (0.56)	=	ND (0.39)	=	ND (0.39)	=	ND (0.39)	ND (0.39)	ND (0.39)	=	1.4 J
2-Butanone (MEK)	300	ND (3.2)	-	ND (3.2)	-	ND (3.2)	-	ND (2.5)	-	ND (2.5)	-	10.2	ND (2.5)	ND (2.5)	=	23.2 J
Carbon disulfide	700	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)	-	2.4 J
Carbon tetrachloride Chlorobenzene	1 50	ND (0.23) ND (0.35)	-	ND (0.23) ND (0.35)	-	ND (0.23) ND (0.35)	-	ND (0.24) ND (0.27)	-	ND (0.24) ND (0.27)	-	ND (0.24) ND (0.27)	ND (0.24) ND (0.27)	ND (0.24) ND (0.27)	-	0.73 J ND (0.68)
Chloroethane	50	ND (0.35) ND (0.39)		ND (0.35) ND (0.39)	-	ND (0.35) ND (0.39)	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)		ND (0.27) ND (0.56)	ND (0.27) ND (0.56)	ND (0.27) ND (0.56)	-	ND (0.68)
Chloroform	70	ND (0.25)	-	ND (0.25)	-	0.47 J	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	ND (0.20)	ND (0.20)	-	4.1
Chloromethane	-	ND (0.36)	-	ND (0.36)	-	ND (0.36)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (0.83)
Cyclohexane	-	ND (0.18)	-	ND (0.18)	-	ND (0.18)	-	ND (0.37)	=	ND (0.37)	-	ND (0.37)	ND (0.37)	ND (0.37)	=	ND (0.93)
1,2-Dibromo-3-chloropropane	0.02	ND (1.3)	-	ND (1.3)	-	ND (1.3)		ND (1.2)	-	ND (1.2)		ND (1.2)	ND (1.2)	ND (1.2)		ND (2.9)
Dibromochloromethane	1	ND (0.19)	-	ND (0.19)	-	ND (0.19)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	-	ND (0.62)
1,2-Dibromoethane	0.03	ND (0.16)	-	ND (0.16)	-	ND (0.16)	=	ND (0.23)	=	ND (0.23)	-	ND (0.23)	ND (0.23)	ND (0.23)	=	ND (0.57)
1,2-Dichlorobenzene 1,3-Dichlorobenzene	600 600	ND (0.20) ND (0.31)	<del>                                     </del>	ND (0.20) ND (0.31)	-	ND (0.20) ND (0.31)	<del>  </del>	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)		ND (0.16) ND (0.26)	ND (0.16) ND (0.26)	ND (0.16) ND (0.26)	-	1.1 J ND (0.64)
1,4-Dichlorobenzene	75	ND (0.31)	<del>                                     </del>	ND (0.31)	-	ND (0.31)	-	ND (0.24)	-	ND (0.24)		ND (0.24)	ND (0.24)	ND (0.24)	-	ND (0.59)
Dichlorodifluoromethane	1000	ND (0.63)	=	ND (0.63)	-	ND (0.63)	-	ND (0.73)	=	ND (0.73)	-	ND (0.73)	ND (0.73)	ND (0.73)	=	ND (1.8)
1,1-Dichloroethane	50	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	-	ND (0.87)
1,2-Dichloroethane	2	ND (0.22)	-	ND (0.22)	-	3.3	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	ND (0.30)	2.6	-	57.8
1,1-Dichloroethene	1	ND (0.34)	-	ND (0.34)	-	ND (0.34)	=	ND (0.50)	÷	ND (0.50)	-	ND (0.50)	ND (0.50)	ND (0.50)	÷	ND (1.2)
cis-1,2-Dichloroethene	70	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	ND (0.33)	-	ND (0.82)
trans-1,2-Dichloroethene	100	ND (0.38) ND (0.28)	-	ND (0.38) ND (0.28)	-	ND (0.38)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	ND (0.51) ND (0.43)	ND (0.51) ND (0.43)	-	ND (1.3) ND (1.1)
1,2-Dichloropropane cis-1,3-Dichloropropene	1	ND (0.28) ND (0.15)	-	ND (0.28) ND (0.15)	-	ND (0.28) ND (0.15)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)		ND (0.43) ND (0.28)	ND (0.43) ND (0.28)	ND (0.43) ND (0.28)	-	ND (1.1) ND (0.71)
trans-1,3-Dichloropropene	-	ND (0.21)	- 1	ND (0.21)	-	ND (0.21)	-	ND (0.32)	_	ND (0.32)	-	ND (0.32)	ND (0.32)	ND (0.32)	-	ND (0.79)
1,4-Dioxane	-	ND (73)	-	ND (73)	-	ND (73)	-	ND (51)	-	ND (51)	-	ND (51)	ND (51)	ND (51)	-	ND (130)
Ethylbenzene	700	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.40)	÷	ND (0.40)	-	ND (0.40)	ND (0.40)	ND (0.40)	-	ND (0.99)
Freon 113	-	ND (0.77)	-	ND (0.77)	-	ND (0.77)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	ND (0.45)	ND (0.45)	-	ND (1.1)
2-Hexanone	-	ND (1.7)	- +	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	ND (1.7)	ND (1.7)	-	ND (4.3)
Isopropylbenzene Methyl Acetate	700 7000	ND (0.22) ND (1.5)	-	ND (0.22) ND (1.5)	-	ND (0.22) ND (1.5)	-	ND (0.26) ND (3.1)	-	ND (0.26) ND (3.1)	-	ND (0.26) ND (3.1)	ND (0.26) ND (3.1)	ND (0.26) ND (3.1)	-	ND (0.64) ND (7.7)
Methylcyclohexane	7000	ND (1.5) ND (0.15)	-	ND (1.5)	-	ND (1.5)	-	ND (3.1)	-	ND (3.1)		ND (5.1)	ND (3.1)	ND (3.1) ND (0.22)	-	ND (7.7) ND (0.54)
Methyl Tert Butyl Ether	70	ND (0.29)	-	ND (0.29)	-	ND (0.29)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	-	ND (0.66)
4-Methyl-2-pentanone(MIBK)	-	ND (1.5)	-	ND (1.5)	-	ND (1.5)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	ND (1.1)	ND (1.1)	-	ND (2.6)
Methylene chloride Styrene	3 100	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.86) ND (0.30)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	ND (0.81) ND (0.26)	ND (0.81) ND (0.26)	-	ND (2.0) ND (0.64)
1,1,2,2-Tetrachloroethane	100	ND (0.20)	-	ND (0.30)	-	ND (0.30)	-	ND (0.39)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	ND (0.26)	-	92.4
Tetrachloroethene	1	ND (0.25)	-	ND (0.25)	-	ND (0.25)	- 1	ND (0.35)	-	ND (0.35)	-	ND (0.35)	ND (0.35)	ND (0.35)	-	7.9
Toluene	600	ND (0.44)		ND (0.44)	-	ND (0.44)	-	ND (0.22)	-	ND (0.22)		ND (0.22)	ND (0.22)	ND (0.22)	-	ND (0.55)
1,2,3-Trichlorobenzene 1,2,4-Trichlorobenzene	9	ND (0.24) ND (0.22)	-	ND (0.24) ND (0.22)	-	ND (0.24) ND (0.22)	-	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	-	ND (0.66) ND (0.56)
1,1,1-Trichloroethane	30	ND (0.25)	-	ND (0.25)	-	ND (0.25)	- 1	ND (0.32)	- 1	ND (0.32)	- +	ND (0.32)	ND (0.32)	ND (0.32)	-	ND (0.80)
1,1,2-Trichloroethane	3	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	ND (0.28)	-	0.81 J
Trichloroethene	1	ND (0.50)	-	ND (0.50)	-	1.4 ND (0.33)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	ND (0.25)	ND (0.25)	-	31.3 ND (0.70)
Trichlorofluoromethane Vinyl chloride	2000	ND (0.33) ND (0.41)	<del>                                     </del>	ND (0.33) ND (0.41)	-	ND (0.33) ND (0.41)		ND (0.28) ND (0.17)	-	ND (0.28) ND (0.17)		ND (0.28) ND (0.17)	ND (0.28) ND (0.17)	ND (0.28) ND (0.17)	-	ND (0.70) ND (0.44)
m,p-Xylene	-	ND (0.41)		ND (0.40)		ND (0.40)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	ND (0.45)	ND (0.45)	-	ND (1.1)
o-Xylene	-	ND (0.19)	-	ND (0.19)	-	ND (0.19)		ND (0.20)	-	ND (0.20)		ND (0.20)	ND (0.20)	ND (0.20)	-	ND (0.50)
Xylene (total) Total VOCs	1000	ND (0.19) <b>0</b>	<del>                                     </del>	ND (0.19) 0	-	ND (0.19) 23.27		ND (0.20) 0	-	ND (0.20) <b>0</b>	-	ND (0.20) 10.2	ND (0.20) <b>0</b>	ND (0.20) 2.6	-	ND (0.50) <b>514.78</b>
Total VOC3			<u> </u>		1	23.21	1					10.2		2.0		314./0
GC/MS Volatile TIC																
Total TIC, Volatile	-	0	- 1	0	-	0		0			-		0		J -	0
Total Alkanes	-	0	-	0	-	0	-	0	-	0	-	0	0	0	-	0
Metals Analysis																
Chromium	70	<10	<10	<10	<10	1390 a	418	<10	<10	<10	<10	- 1	- 1	<10	<10	718
Iron	300	-	<100	-		-	9150	-	<100	-	<100		-	-	<100	-
Sodium	50000	103000	-	93100	-	482000 a		<10000		<10000	-	-		89300		1060000
General Chemistry																
Solids, Total Dissolved	500000	677000	1   -	683000	-	2730000	- 1	<10000	-	<10000			1 - 1	452000	1 - 1	3860000
Sulfate	250000		-	278000		350000		<10000	-		-	-	-	181000	-	1880000
																•

## 1971 1971 1971 1971 1971 1971 1971 19	Sample ID		1RND2 ISCO-MW2	1RND2 ISCO-MW3	1RND2 ISCO-MW3	1RND2 ISCO-MW4	1RND2 ISCO-MW4	1RND2 ISCO-MW5	1RND2 ISCO-MW5	1RND2 ISCO-MW6	1RND2 ISCO-MW6	1RND2 DUP	1RND2 DUP	1RND2 ISCO-MW7	1RND2 ISCO-MW7	1RND2 ISCO-MW8
Company		NJ CLASS IIA														JB70619-12
The control of the co		GROUNDWATER QUALITY	7/1/2014	7/1/2014	7/1/2014	6/30/2014	6/30/2014	6/30/2014	6/30/2014	7/1/2014	7/1/2014	7/1/2014	7/1/2014	6/30/2014	6/30/2014	7/1/2014
Column		CRITERIA (7/22/2010)	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW - FILTERED	GW	GW-FILTERED	GW
1	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Column	Volatile Organic Compounds (VOCs)															
Company   Comp	Acetone	6000	-		-	ND (2.6)	-		-		-		-		-	36.8
Secretary   1		1	-		-		-		-		-		-		-	ND (0.21)
Second   1		-	-	` '	-		-		-		-		-		-	ND (0.49)
March			-		-	` '	-	· · · · · · · · · · · · · · · · · · ·	-	` '	-	` '			-	ND (0.28)
Company   12			-	` '	-	` '	-	(/	=	` '	-	` '	-		-	ND (0.31)
Something 150			-		-		-	· · · · · · · · · · · · · · · · · · ·	-		-	· · ·			-	ND (0.39) ND (2.5)
Contribution   1					-		-		-		-		+ -		_	ND (0.50)
Company   Comp			_	` '	_		_	` '	_		_		_	, ,	1 -	0.35 J
Control   Cont			-	` '	-	` '	-	` '	-	` '	-	· · ·	- 1			ND (0.27)
Company	Chloroethane	-	-	` '	-	` '	-	` '	-	` '	-	` '	-	, ,	-	ND (0.56)
Company   Comp	Chloroform	70	-	ND (0.20)	-		-	ND (0.20)	-	ND (0.20)	-		-	1.6	-	1.4
March   Marc	Chloromethane	-	-	ND (0.33)		ND (0.33)	-	ND (0.33)		ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Manufacturing   1		-	-		-	` '	-		-	· · · · · · · · · · · · · · · · · · ·	- 1	· · ·	- 1		-	ND (0.37)
30   10   10   10   10   10   10   10			-		-			<del></del>	-	· · · · · · · · · · · · · · · · · · ·			1		-	ND (1.2)
13C9-1200-1200-1200-1200-1200-1200-1200-120			=		-		-	· · · · · · · · · · · · · · · · · · ·	-		-	· · ·			-	ND (0.25)
1300-1500-1500-1500-1500-1500-1500-1500-			-		-		-		-		-		+ - +		-	ND (0.23)
Control   Cont			=	` '	-		-		-		-		+ +		-	ND (0.16)
Content   Cont			-	` '	1	` '	-	` '	<del>                                     </del>	` '	-	` '	+ - +	, ,	-	ND (0.26) ND (0.24)
10.00000000000000000000000000000000000			-		-	<del></del>		· · · · · · · · · · · · · · · · · · ·	-	· ' '	-	· · ·	+ +		-	ND (0.24) ND (0.73)
Calcalementary   Calc			-	` '	-	` '	-		-		_		+	, ,	-	ND (0.75) ND (0.35)
Continue			-	` '	-	` '	-		-	` '	-	` '	-	, ,	-	0.48 J
Section   Sect			-	` '	-		-		-		-		-		-	ND (0.50)
1	cis-1,2-Dichloroethene	70	-	57.2	-	ND (0.33)	-	4.3	-	4.7	-	4.8	-	ND (0.33)	-	0.61 J
Exp. Continue proper	trans-1,2-Dichloroethene	100	=	0.57 J	-	ND (0.51)	=	ND (0.51)	=	ND (0.51)	=	ND (0.51)	=	ND (0.51)	ē.	ND (0.51)
Trans La Diffusionegone	1,2-Dichloropropane	1	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)
1.4 Decided   1.5 Decided		=	-	` '	-		-	` '	-		-		-		-	ND (0.28)
Targetonises   720	/		-	` '	-	` '	-	· · · · · · · · · · · · · · · · · · ·	-	` '	-	· · ·	-		-	ND (0.32)
Fig. 1			-		-	` '	-	1- /	-		-	· · · · ·			-	ND (51)
2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-2-	'	700	-	` '	-	` '	-		-	` '	-	` '		(/	-	ND (0.40)
Manual Control		-		` '	-	` '		` '	-	· · · · · · · · · · · · · · · · · · ·	-	` '	<del>-</del> -		-	ND (0.45) ND (1.7)
Mode		700	_		_		_		_		_		_		_	ND (0.26)
Motiny for short detection of the short of t			-		-	<del></del>	-		-	· · · · · · · · · · · · · · · · · · ·	-	· · ·	-			ND (3.1)
Steffely 2 pertamen(MISS)   NO (13)   NO (13	Methylcyclohexane	-	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
Morphisen sharing  1		70	-		-		-		-		-		-		-	ND (0.26)
Syrene   100		-	-		-		-		-		-		-		-	ND (1.1)
11.2.2-Tetrathrocethane			-		-		-		-		-	· · ·	+	1 /	-	ND (0.81) ND (0.26)
Transference   1			-		-		-		-		-		-		-	ND (0.39)
12.3 Trichiorobenseme		1	-		-	` '	-		-		-		J -		J -	0.58 J
1.2.4-Trichrorebenene		600	-		-		-		-		-				-	ND (0.22)
1.1.1-frickhorechane   30		- 0	=		-	` '	-		-		-	· · ·	+ - +	1 /	_	ND (0.26) ND (0.22)
1			-		-	` '	-		-	· ' '	-	· · ·	+ +			ND (0.22) ND (0.32)
Trichforfourcemethane			<u>-</u>		-		<u>-</u>		<u>-</u>		<u>-</u>				-	ND (0.28)
Vary chloride 1 - ND (0.17) -			-		-		-		-		-		-		-	2.9
Imp-lykene			-		-		-				-		+ - +		-	ND (0.28)
0-Xylene		1	-		-	` '	-		-	· ' '	-	· · ·	+	1 /	-	ND (0.17) ND (0.45)
No   No   No   No   No   No   No   No		-	-		<del>'</del>		-		-		-		+ +		-	ND (0.43)
Common   C	Xylene (total)	1000	<u></u>	ND (0.20)	<u> </u>	ND (0.20)	<u> </u>	ND (0.20)	<u> </u>	ND (0.20)	<u> </u>	ND (0.20)	<u> </u>	ND (0.20)		ND (0.20)
Total TIC, Volatile 0 - 0 - 17 J - 0 - 0 - 0 - 0 - 17 J - 0 - 0 - 0 - 1 - 1 - 1 - 1 - 1 - 1 - 1	Total VOCs	-		150.37		2.61		48.25		15.99		16.19		29.76		43.12
Total TIC, Volatile	CC/MS Volotile TIC															
Total Alkanes 0 0 - 0 - 0 - 0 - 0 - 0 - 0 - 0 -		_		1 0 1	1		1 -	17 11	1 -	1 0	1 -		1	1 0		
Metals Analysis  Chromium 70 769 4480 5320 13.1 < 10 < 10 27.5 < 10 25.3 < 10 490 473   Iron 300 < 100   -   1800   -   1800   -   1800   -   5760   Sodium 5000 -   211000   -   90000   -   153000   -   50000   -   398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -     398000   -																0
Chromium   70   769   4480   5320   13.1   <10   <10   <10   27.5   <10   25.3   <10   490   473		<u>l</u>		<u> </u>	1	ı	1	. · ·	1	· · · · · ·	<u>,                                      </u>	<u> </u>	·		<u> </u>	
Iron 300 <100 - <100 - <100 - 186 - <100 - <100 - 5760   Sodium 50000 - 2110000 - 11800 - 22900 - 46200 - 46600 - 598000 -   General Chemistry Solids, Total Dissolved 50000 - 9030000 - 9030000 - 153000 - 50000 - 3880000 - 3880000 - 3880000 - 588000000 - 5880000 - 5880000000000	Metals Analysis															
Sodium 50000 - 2110000 - 11800 - 12900 - 46200 - 46600 - 598000 - 6  General Chemistry Solids, Total Dissolved 50000 - 9030000 - 9030000 - 153000 - 50000 - 3880000 - 1				4480		13.1		<10		27.5		25.3		490		2820
General Chemistry Solids, Total Dissolved 50000 - 9030000 - 900000 - 153000 - 3980000 - 398000 - 398000 - 398000 - 3050000 - 305000 - 3050000 - 3050000 - 3050000 - 3050000 - 30500000 - 30500000 - 30500000 - 30500000 - 30500000 - 30500000 - 305000								20077								-
Solids, Total Dissolved 500000 - 9030000 - 900000 - 153000 - 50000 - 380000 - 380000 - 380000 - 50000 - 50000 - 50000 - 500000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 500000 - 500000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 500000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50	Soainm	50000	=	2110000	=	11800	-	22900	=	46200	-	46600	1 - 1	598000	-	1760000
Solids, Total Dissolved 500000 - 9030000 - 900000 - 153000 - 50000 - 380000 - 380000 - 380000 - 50000 - 50000 - 50000 - 500000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 500000 - 500000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 500000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50000 - 50	General Chemistry															
	7	500000	- I	9030000	- 1	90000	- 1	153000	- 1	50000	- 1	305000	1 - 1	3980000	- 1	6010000
Sulfate 250000 - 2450000 - 41100 - 74800 - 145000 - 148000 - 703000 - 703000	Sulfate	250000			-	41100	-	74800	-	145000		148000	- 1		-	1910000

Sample ID	NJ CLASS IIA	1RND2_ISCO-MW8 JB70619-12F	1RND2_ISCO-MW9 JB70619-7	1RND2_ISCO-MW9 JB70619-7F	1RND2_IW2-BT2 JB70619-15	1RND2_IW2-BT2 JB70619-15F	1RND2_MW-10S JB70619-1	1RND2_MW-10S JB70619-1F	1RND2_MW-14S-D JB70619-13	1RND2_MW-14S-D JB70619-13F	1RND2_MW-14S-S JB70619-14	1RND2_MW-14S-S JB70619-14F	1RND2_PZ-1S JB70619-8	1RND2_PZ-1S JB70619-8F	1RND3_FB _080714 JB73631-17
Lab Sample ID Sample Date	GROUNDWATER QUALITY CRITERIA (7/22/2010)	7/1/2014	7/1/2014	7/1/2014	7/1/2014	7/1/2014	6/30/2014	6/30/2014	7/1/2014	7/1/2014	7/1/2014	7/1/2014	7/1/2014	7/1/2014	8/7/2014
Matrix	ug/L	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	WATER
Unit Volatile Organic Compounds (VOCs)		ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Acetone	6000	-	5.0	J -	9.2	J -	ND (2.6)	-	ND (2.6)	-	ND (2.6)	-	3.7	J -	ND (2.6)
Benzene Bromochloromethane	1	-	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	=	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	=	ND (0.21) ND (0.49)
Bromodichloromethane	1	-	ND (0.49)	-	ND (0.28)	=	ND (0.28)	-	ND (0.49)	-	ND (0.28)	-	ND (0.28)	=	ND (0.49)
Bromoform	4	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)
Bromomethane 2-Butanone (MEK)	10 300	-	ND (0.39) ND (2.5)	-	ND (0.39) <b>2.8</b>	- 1 -	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	=	ND (0.39) ND (2.5)
Carbon disulfide	700	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)
Carbon tetrachloride	1	-	ND (0.24)	-	ND (0.24)	=	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	=	ND (0.24)
Chlorobenzene Chloroethane	50 -	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)	-	ND (0.27) ND (0.56)
Chloroform	70	-	ND (0.20)	=	ND (0.20)	=	ND (0.20)	-	ND (0.20)	-	ND (0.20)	=	0.27	J -	ND (0.20)
Chloromethane	-	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Cyclohexane 1,2-Dibromo-3-chloropropane	0.02	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)		ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)
Dibromochloromethane	1	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)
1,2-Dibromoethane 1,2-Dichlorobenzene	0.03 600	-	ND (0.23) ND (0.16)		ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)
1,3-Dichlorobenzene	600	-	ND (0.16)	-	ND (0.26)	=	ND (0.26)	-	ND (0.16)	-	ND (0.26)	-	ND (0.26)	=	ND (0.26)
1,4-Dichlorobenzene	75	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)
Dichlorodifluoromethane 1,1-Dichloroethane	1000 50	-	ND (0.73) ND (0.35)	-	ND (0.73) ND (0.35)	-	ND (0.73) ND (0.35)	-	ND (0.73) ND (0.35)	-	ND (0.73) ND (0.35)	-	ND (0.73) ND (0.35)	=	ND (0.73) ND (0.35)
1,2-Dichloroethane	2	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	3.8	-	ND (0.30)
1,1-Dichloroethene	1	-	ND (0.50)	-	ND (0.50)	=	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	=	ND (0.50)
cis-1,2-Dichloroethene trans-1,2-Dichloroethene	70 100	-	<b>0.43</b> ND (0.51)	J -	12.9 2.2	-	ND (0.33) ND (0.51)	-	ND (0.33) ND (0.51)	-	ND (0.33) ND (0.51)	-	ND (0.33) ND (0.51)	-	ND (0.33) ND (0.51)
1,2-Dichloropropane	1	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)
cis-1,3-Dichloropropene	-	-	ND (0.28)	-	ND (0.28)	=	ND (0.28)	-	ND (0.28)	-	ND (0.28)	=	ND (0.28)	=	ND (0.28)
trans-1,3-Dichloropropene 1.4-Dioxane	-	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)
Ethylbenzene	700	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)
Freon 113	-	-	ND (0.45)	-	ND (0.45)	-	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)	-	ND (0.45)	-	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)
2-Hexanone Isopropylbenzene	700	-	ND (1.7) ND (0.26)	-	ND (1.7) ND (0.26)	-	ND (1.7) ND (0.26)	-	ND (1.7)	-	ND (1.7) ND (0.26)	-	ND (1.7) ND (0.26)	-	ND (1.7)
Methyl Acetate	7000	-	ND (3.1)	-	ND (3.1)	-	ND (3.1)	-	ND (3.1)	-	ND (3.1)	-	ND (3.1)	-	ND (3.1)
Methylcyclohexane Methyl Tert Butyl Ether	- 70	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)
4-Methyl-2-pentanone(MIBK)	-	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)
Methylene chloride Styrene	3 100	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)
1,1,2,2-Tetrachloroethane	1	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)
Tetrachloroethene Toluene	1 600	-	0.83 ND (0.22)	J -	<b>0.76</b> ND (0.22)	J -	ND (0.35) ND (0.22)	-	ND (0.35) ND (0.22)	-	ND (0.35) ND (0.22)	-	ND (0.35) ND (0.22)	-	ND (0.35) ND (0.22)
1,2,3-Trichlorobenzene	-	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
1,2,4-Trichlorobenzene 1,1,1-Trichloroethane	9 30	-	ND (0.22) ND (0.32)		ND (0.22) ND (0.32)		ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)
1,1,2-Trichloroethane	3	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
Trichloroethene Trichlorofluoromethane	1 2000	-	15.8 ND (0.28)		<b>26.8</b> ND (0.28)		ND (0.25) ND (0.28)	-	ND (0.25) ND (0.28)	-	ND (0.25) ND (0.28)		1.7 ND (0.28)		ND (0.25) ND (0.28)
Vinyl chloride	1	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)
m,p-Xylene o-Xylene	-	-	ND (0.45) ND (0.20)		ND (0.45) ND (0.20)	= =	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)		ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)
Xylene (total) Total VOCs	1000	-	ND (0.20) 22.66	-	ND (0.20) <b>54.66</b>	-	ND (0.20) <b>0</b>	-	ND (0.20) 0	-	ND (0.20) 0	-	ND (0.20) 9.47	-	ND (0.20) <b>0</b>
	<u> </u>		22.00		54.00		U	<u> </u>			U		5.47		
GC/MS Volatile TIC				Ţ											
Total TIC, Volatile Total Alkanes	-	-	0	-	0	-	0	-	0	-	0	-	<b>11</b> 0	J -	0
		<u> </u>	•			•	<u> </u>					•		•	
Metals Analysis Chromium	70	2900	86.9	18.3	584	<10	<10	<10	<10	<10	<10	<10	105	86	<10
Iron	300	<100	-	<100	-	<100	-	<100	-	<100	-	<100	-	1880	-
Sodium	50000	-	147000	-	197000	-	<10000	-	125000	-	111000	-	174000	-	<10000
General Chemistry															
Solids, Total Dissolved Sulfate	500000 250000				1160000 465000		32000 18700		731000 218000		669000 195000		1040000 128000		<10000 <10000
Sanate	250000		212000		403000		10/00	1	210000	1	193000	-	120000		10000

	Sample ID	1RND3_FB_080714	1RND3_FB_08081	4 1RND3_FB_080814	ТВ	1RND3 ISCO-MW1	1RND3 ISCO-MW1	1RND3 ISCO-MW1 ASC 0808	14 1RND3 ISCO-MW2	1RND3 ISCO-MW2	1RND3 ISCO-MW2 ASC 08083	L4 1RND3 ISCO-MW3	1RND3 ISCO-MW3	1RND3 ISCO-MW4	1RND3 ISCO-MW4
	Lah Sample ID	JB73631-17F	JB73631-16	JB73631-16F	JB73631-18	JB73631-11	JB73631-11F	JB73631-20	JB73631-13	JB73631-13F	JB73631-22	JB73631-6	JB73631-6F	JB73631-5	JB73631-5F
	Sample Date GROUNDWATER QUALITY CRITERIA (7/22/2010)	8/7/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/7/2014	8/7/2014	8/7/2014	8/7/2014
	Matrix ug/L	WATER-FILTERED	WATER	WATER-FILTERED	WATER	GW	GW-FILTERED	GW	GW	GW-FILTERED	GW	GW	GW-FILTERED	GW	GW-FILTERED
Valatila Organia Compounds (V	Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (V	6000	_	ND (2.6)		ND (2.6)	ND (2.6)	_	ND (2.6)	375		302	47.3		ND (2.6)	
Acetone Benzene	1	-	ND (2.6)	-	ND (0.21)	ND (0.21)	-	ND (2.0) ND (0.21)	2.3	-	3.4	ND (0.21)	-	ND (0.21)	<del></del>
Bromochloromethane	=	-	ND (0.49)	-	ND (0.49)	ND (0.49)	-	ND (0.49)	ND (0.49)	-	ND (0.49)	ND (0.49)	-	ND (0.49)	-
Bromodichloromethane	1	-	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	-
Bromoform	4	-	ND (0.31)	-	ND (0.31)	ND (0.31)	-	ND (0.31)	ND (0.31)	-	ND (0.31)	ND (0.31)	-	ND (0.31)	-
Bromomethane	10 300	=	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	ND (0.39) ND (2.5)	=	ND (0.39) ND (2.5)	41.2 316	-	70.8 357	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-
2-Butanone (MEK) Carbon disulfide	700	-	ND (0.50)	-	ND (0.50)	ND (2.5) ND (0.50)	-	ND (2.5)	38	-	25.8	ND (0.50)	-	ND (0.50)	<del>-</del>
Carbon tetrachloride	1	-	ND (0.24)	-	ND (0.24)	ND (0.24)	-	ND (0.24)	0.27 J	-	ND (0.24)	ND (0.24)	-	ND (0.24)	-
Chlorobenzene	50	-	ND (0.27)	-	ND (0.27)	ND (0.27)	-	ND (0.27)	ND (0.27)	-	ND (0.27)	ND (0.27)	-	ND (0.27)	-
Chloroethane	-	-	ND (0.56)	-	ND (0.56)	ND (0.56)	-	ND (0.56)	ND (0.56)	-		J ND (0.56)	-	ND (0.56)	<del>-</del>
Chloroform Chloromethane	70	-	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	4.2	-	3.6 4.2	ND (0.20) ND (0.33)	-	ND (0.20) ND (0.33)	-
Cyclohexane		-	ND (0.37)	-	ND (0.37)	ND (0.33)	-	ND (0.37)	ND (0.37)	-	ND (0.37)	ND (0.37)	-	ND (0.37)	-
1,2-Dibromo-3-chloropropane	0.02	-	ND (1.2)	-	ND (1.2)	ND (1.2)	-	ND (1.2)	ND (1.2)	-	ND (1.2)	ND (1.2)	<u>-</u>	ND (1.2)	-
Dibromochloromethane	1	-	ND (0.25)	-	ND (0.25)	ND (0.25)	-	ND (0.25)	ND (0.25)	-	ND (0.25)	ND (0.25)	-	ND (0.25)	-
1,2-Dibromoethane	0.03	-	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	<del>-</del>
1,2-Dichlorobenzene 1,3-Dichlorobenzene	600 600	-	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	0.52 J ND (0.26)	-	5.1 ND (0.26)	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	-
1,4-Dichlorobenzene	75	-	ND (0.24)	-	ND (0.24)	ND (0.26) ND (0.24)	-	ND (0.26) ND (0.24)	ND (0.26) ND (0.24)	<del>                                     </del>	, ,	J ND (0.24)	<del>-</del> -	ND (0.24)	-
Dichlorodifluoromethane	1000	-	ND (0.73)	-	ND (0.73)	ND (0.73)	-	ND (0.73)	ND (0.73)	<u> </u>	ND (0.73)	ND (0.73)		ND (0.73)	-
1,1-Dichloroethane	50	=	ND (0.35)	-	ND (0.35)	ND (0.35)	=	ND (0.35)	ND (0.35)	-	ND (0.35)	ND (0.35)	-	ND (0.35)	-
1,2-Dichloroethane	2	-	ND (0.30)	-	ND (0.30)	2.9	-	2.7	485	-	554	ND (0.30)	-	ND (0.30)	-
1,1-Dichloroethene cis-1,2-Dichloroethene	70	-	ND (0.50) ND (0.33)	-	ND (0.50) ND (0.33)	ND (0.50) 1.1	-	ND (0.50) ND (0.33)	ND (0.50) <b>0.63</b> J	-	ND (0.50) 0.34	ND (0.50) J 87.1	-	ND (0.50) ND (0.33)	-
trans-1,2-Dichloroethene	100	-	ND (0.51)	-	ND (0.51)	19.3	-	15.1	ND (0.51)	-	ND (0.51)	27.1	-	30.3	<del>                                     </del>
1,2-Dichloropropane	1	-	ND (0.43)	-	ND (0.43)	ND (0.43)	-	ND (0.43)	ND (0.43)	-	ND (0.43)	ND (0.43)	-	ND (0.43)	-
cis-1,3-Dichloropropene	-	=	ND (0.28)	-	ND (0.28)	ND (0.28)	=	ND (0.28)	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	-
trans-1,3-Dichloropropene	=	=	ND (0.32)	-	ND (0.32)	ND (0.32)	=	ND (0.32)	ND (0.32)	-	ND (0.32)	ND (0.32)	-	ND (0.32)	-
1,4-Dioxane	700	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	ND (51) ND (0.40)	-	ND (51) ND (0.40)	ND (51)	-	ND (51) ND (0.40)	ND (51)	-	ND (51) ND (0.40)	-
Ethylbenzene Freon 113	700	-	ND (0.45)	-	ND (0.40) ND (0.45)	ND (0.40) ND (0.45)	-	ND (0.40) ND (0.45)	ND (0.40) ND (0.45)	-	ND (0.40) ND (0.45)	ND (0.40) ND (0.45)	-	ND (0.45)	-
2-Hexanone	-	=	ND (1.7)	-	ND (1.7)	ND (1.7)	=	ND (1.7)	ND (1.7)	-	ND (1.7)	ND (1.7)	-	ND (1.7)	=
Isopropylbenzene	700	=	ND (0.26)	-	ND (0.26)	ND (0.26)	=	ND (0.26)	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
Methyl Acetate	7000	-	ND (3.1)	-	ND (3.1)	ND (3.1)	-	ND (3.1)	ND (3.1)	-	ND (3.1)	ND (3.1)	-	ND (3.1)	-
Methylcyclohexane Methyl Tert Butyl Ether	70	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-
4-Methyl-2-pentanone(MIBK)	-	-	ND (1.1)	-	ND (1.1)	ND (1.1)	-	ND (1.1)	ND (1.1)	-	ND (1.1)	ND (1.1)	-	ND (1.1)	-
Methylene chloride	3	-	ND (0.81)	-	ND (0.81)	ND (0.81)	-	ND (0.81)	3.4	-	3.3	ND (0.81)	-	ND (0.81)	-
Styrene 1,1,2,2-Tetrachloroethane	100	-	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	ND (0.26) 17.6	-	ND (0.26) 16.4	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-
Tetrachloroethene	1		ND (0.35)		ND (0.35)	ND (0.35)		ND (0.35)	3.7	<u> </u>	1.6	2		ND (0.35)	-
Toluene	600	=	ND (0.22)	-	ND (0.22)	ND (0.22)	=	ND (0.22)	0.6 J		ND (0.22)	ND (0.22)	-	ND (0.22)	-
1,2,3-Trichlorobenzene 1,2,4-Trichlorobenzene	9	-	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	ND (0.26) <b>0.6</b> J	-	1.6 13.5	J ND (0.26) ND (0.22)	+ -	ND (0.26) ND (0.22)	-
1,1,1-Trichloroethane	30	-	ND (0.32)		ND (0.32)	ND (0.32)	-	ND (0.32)	ND (0.32)	-	ND (0.32)	ND (0.32)	<u>-</u>	ND (0.32)	-
1,1,2-Trichloroethane	3	-	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	0.46 J	-	ND (0.28)	ND (0.28)	-	ND (0.28)	
Trichloroethene Trichlorofluoromethane	2000	-	ND (0.25) ND (0.28)	-	ND (0.25) ND (0.28)	ND (0.25) ND (0.28)	-	ND (0.25) ND (0.28)	16.3 ND (0.28)	<del>  -</del>	6.5 ND (0.28)	71.5 ND (0.28)	-	0.64 ND (0.28)	J -
Vinyl chloride	1		ND (0.17)		ND (0.17)	ND (0.17)		ND (0.17)	ND (0.17)	<u> </u>	ND (0.17)	ND (0.17)	<u>-</u> -	ND (0.17)	-
m,p-Xylene	-	-	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	
o-Xylene Xylene (total)	1000	-	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	ND (0.20) ND (0.20)	+ -	ND (0.20) ND (0.20)	<del></del>
Total VOCs	-		0		0	23.3		17.8	1310.08			235		30.94	
CO (NAC V. L. III - TIC															
GC/MS Volatile TIC Total TIC, Volatile			0	- 1	0 1	0		0	35.3 J		37.7	J 19	J] - [	0	
Total Alkanes		-			0	0	-	0	0 J			0			-
											·				
Metals Analysis		150				1			4000	762	, ·	2000	F200		
Chromium Iron	70 300	<10 <100	<10	<10 <100	-	<10	<10 128	-	1090	<b>762</b> <100	-	3450	<b>5280</b> <100	<10	<10 <100
Sodium	50000		<10000		-	50400		-	2050000		-	1110000	-	14100	-
General Chemistry Solids, Total Dissolved	500000		<10000			262000			4850000	1	1	5330000		84000	
Sulfate	250000	-	<10000	-	-		-	-	2630000	-		1790000	-	45000	-
	•				•	•		•	•		•			•	

Sample ID	NJ CLASS IIA	1RND3_ISCO-MW5	1RND3_ISCO-MW5	1RND3_ISCO-MW5 ASC _080814		1RND3_ISCO-MW6	1RND3_ISCO-MW7	1RND3_ISCO-MW7	1RND3-DUP_080714			1RND3_ISCO-MW8	1RND3_ISCO-MW9 JB73631-10	1RND3_ISCO-MW9 JB73631-10F	1RND3_ISCO-MW9 ASC _080814
Lab Sample ID Sample Date	GROUNDWATER QUALITY	JB73631-12 8/8/2014	JB73631-12F 8/8/2014	JB73631-21 8/8/2014	JB73631-1 8/7/2014	JB73631-1F 8/7/2014	JB73631-4 8/7/2014	JB73631-4F 8/7/2014	JB73631-9 8/7/2014	JB73631-9F 8/7/2014	JB73631-2 8/7/2014	JB73631-2F 8/7/2014	8/8/2014	8/8/2014	JB73631-19 8/8/2014
Matrix	CRITERIA (7/22/2010)	GW	GW-FILTERED	GW	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW
Unit	ug/L it	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs)															
Acetone	6000	ND (2.6)	-	ND (2.6)	ND (2.6)	-	28.4	-	29	-	39.2	-	ND (2.6)	-	ND (2.6)
Benzene	1	ND (0.21)	-	ND (0.21)	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Bromochloromethane Bromodichloromethane	1	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)
Bromoform	4	ND (0.28) ND (0.31)	-	ND (0.28)	ND (0.28) ND (0.31)	-	ND (0.28) ND (0.31)	-	ND (0.28)	-	ND (0.28) ND (0.31)	-	ND (0.28) ND (0.31)	-	ND (0.28) ND (0.31)
Bromomethane	10	ND (0.31)		ND (0.39)	ND (0.31)	-	2.1	_	2	-	ND (0.31)		ND (0.31)	_	ND (0.39)
2-Butanone (MEK)	300	ND (2.5)	-	ND (2.5)	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)
Carbon disulfide	700	ND (0.50)	-	ND (0.50)	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)
Carbon tetrachloride	1	ND (0.24)	-	ND (0.24)	ND (0.24)	-	0.63	J -	0.68 J	-	0.44 J	-	ND (0.24)	-	ND (0.24)
Chlorobenzene	50	ND (0.27)	-	ND (0.27)	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)
Chloroethane	-	ND (0.56)	-	ND (0.56)	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)
Chloroform	70	ND (0.20)	-	ND (0.20)	ND (0.20)	-	1.6	-	1.6	-	1.6	-	ND (0.20)	-	ND (0.20)
Chloromethane	-	ND (0.33)	-	ND (0.33)	ND (0.33)	-	0.6	J -	0.5 J	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)
Cyclohexane 1,2-Dibromo-3-chloropropane	0.02	0.65 J ND (1.2)	-	ND (0.37) ND (1.2)	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)
1,2-Dibromo-3-chloropropane Dibromochloromethane	0.02	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)
1,2-Dibromoethane	0.03	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	-	ND (0.23)		ND (0.23)		ND (0.23)	-	ND (0.23)
1,2-Dichlorobenzene	600	ND (0.16)	-	ND (0.16)	ND (0.16)	-	ND (0.16)	-	ND (0.16)	- 1	ND (0.16)	-	ND (0.16)	-	ND (0.16)
1,3-Dichlorobenzene	600	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	÷	ND (0.26)
1,4-Dichlorobenzene	75	ND (0.24)	-	ND (0.24)	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)
Dichlorodifluoromethane	1000	ND (0.73)	-	ND (0.73)	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)
1,1-Dichloroethane	50	ND (0.35)	-	ND (0.35)	ND (0.35)	-	ND (0.35)		ND (0.35)		ND (0.35)	-	ND (0.35)		ND (0.35)
1,2-Dichloroethane	2	1.2	-	1	3.4	-	2.5	-	2.5	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)
1,1-Dichloroethene	1	ND (0.50)	-	ND (0.50)	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)
cis-1,2-Dichloroethene	70 100	5.7 ND (0.51)	-	3.1 ND (0.51)	3.7	-	5.9	-	6.8	-	ND (0.33)	-	9.3	-	ND (0.33)
trans-1,2-Dichloroethene 1,2-Dichloropropane	100	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	ND (0.51) ND (0.43)	-	12.3 ND (0.43)	-	11.9 ND (0.43)	-	ND (0.51) ND (0.43)	<del>                                     </del>	19.9 ND (0.43)	-	29.7 ND (0.43)
cis-1,3-Dichloropropene	1	ND (0.43) ND (0.28)	-	ND (0.28)	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.28)
trans-1,3-Dichloropropene	-	ND (0.32)	-	ND (0.32)	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	_	ND (0.32)
1,4-Dioxane	=	ND (51)	-	ND (51)	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)
Ethylbenzene	700	ND (0.40)	-	ND (0.40)	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)
Freon 113	-	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)
2-Hexanone	=	ND (1.7)	÷	ND (1.7)	ND (1.7)	=	ND (1.7)	-	ND (1.7)	-	ND (1.7)	=	ND (1.7)	÷	ND (1.7)
Isopropylbenzene	700	0.79 J	-	0.41 J	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
Methyl Acetate	7000	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1)
Methylcyclohexane Methyl Tert Butyl Ether	70	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)
4-Methyl-2-pentanone(MIBK)	-	ND (1.1)	=	ND (1.1)	ND (1.1)	=	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)		ND (1.1)
Methylene chloride	3	ND (0.81)	-	ND (0.81)	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)
Styrene	100	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
1,1,2,2-Tetrachloroethane Tetrachloroethene	1	ND (0.39) 0.4 J	-	ND (0.39) ND (0.35)	ND (0.39) <b>0.41</b> J	-	ND (0.39) <b>0.66</b>	-	ND (0.39) 0.62 J	-	ND (0.39) <b>0.65</b> J	-	ND (0.39) 1	-	ND (0.39) 0.75 J
Toluene	600	ND (0.22)	-	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,2,3-Trichlorobenzene	-	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
1,2,4-Trichlorobenzene	9	ND (0.22)	-	ND (0.22)	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,1,1-Trichloroethane 1,1,2-Trichloroethane	30	1.4 ND (0.28)	-	0.79 J ND (0.28)	ND (0.32) ND (0.28)	-	ND (0.32) ND (0.28)	-	ND (0.32) ND (0.28)		ND (0.32) ND (0.28)	-	ND (0.32) ND (0.28)	-	ND (0.32) ND (0.28)
Trichloroethene	1	18.4	-	11.2	5.5	-	ND (0.28)	-	2.1	-	2.9	-	14.7	-	11.3
Trichlorofluoromethane	2000	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
Vinyl chloride	1	ND (0.17)	-	ND (0.17)	ND (0.17)	-	ND (0.17)	-	ND (0.17)	-	ND (0.17)	<u> </u>	ND (0.17)	-	ND (0.17)
m,p-Xylene o-Xylene	-	ND (0.45) 1.5	-	ND (0.45) <b>0.91</b> J	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)	-	ND (0.45) ND (0.20)
o-xylene Xylene (total)	1000	1.5	-	0.91 J	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)		ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)
Total VOCs	-	30.04		17.41	13.01		56.69		57.7		44.79		44.9		41.75
GC/MS Volatile TIC								Ţ							
Total TIC, Volatile Total Alkanes	-	5.5 J	-	0	0	-	0	-	0	-	0	-	0	-	0
rotar ARMIES	-	U	-	U	U	-	U	-	U	- 1	U	-	U	-	U
Metals Analysis															
Chromium	70	<10	<10	-	<10	<10	275		294		1950	2020	142	<10	-
Iron	300	-	393	-	-	<100	-	3180	-	5550	-	928	-		-
Sodium	50000	22800	-	-	62700	-	464000	-	505000	-	1330000	-	53600	-	-
General Chemistry															
Solids, Total Dissolved	500000	105000	- 1	-	360000	-	746000	- 1	3720000		5460000		111000	-	<u> </u>
Sulfate	250000	62100	-	-		-			718000	-	1420000		137000		-

Sample II	D	1RND3 IW1-BT2 080714	1 1RND3_IW1-BT2_08071	4 1RND3 MW-10S 0807	714 1RND3 MW-10S 08071	4 1RND3_MW-14SD_080814	1 1RND3 MW-14SD 080814	4 1RND3 MW-14SS 0808	14 1RND3 MW-14SS 0808	314 1RND3_PZ-1S	1RND3_PZ-1S	IRND4_FB _09102014	1 IRND4_FB _09102014	IRND4_FB _09112014	4 IRND4_FB_09112014	IRND4_TB
Lab Sample II		JB73631-3	JB73631-3F	JB73631-8	JB73631-8F	JB73631-15	JB73631-15F	JB73631-14	JB73631-14F	JB73631-7	JB73631-7F	JB76271-11	JB76271-11F	JB76271-21	JB76271-21F	JB76271-12
Sample Dat	e CRITERIA (7/22/2010)	8/7/2014	8/7/2014	8/7/2014	8/7/2014	8/8/2014	8/8/2014	8/8/2014	8/8/2014	8/7/2014	8/7/2014	9/10/2014	9/10/2014	9/11/2014	9/11/2014	9/11/2014
Matri	x ug/L	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	WATER	WATER-FILTERED	WATER	WATER-FILTERED	WATER
Uni Volatile Organic Compounds (VOCs)	it	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Acetone	6000	10.3		ND (2.6)		ND (2.6)		ND (2.6)	_	ND (2.6)	1 - 1	ND (2.6)		ND (2.6)	1 - 1	ND (2.6)
Benzene	1	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)
Bromochloromethane	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)
Bromodichloromethane	1	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)
Bromoform	4	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	=	ND (0.31)	-	ND (0.31)	÷	ND (0.31)	-	ND (0.31)
Bromomethane	10	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)
2-Butanone (MEK)	300	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)	-	ND (2.5)
Carbon disulfide Carbon tetrachloride	700	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)
Chlorobenzene	50	ND (0.24) ND (0.27)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)
Chloroethane	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	_	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)
Chloroform	70	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	1.9	-	3.9	-	ND (0.20)
Chloromethane	=	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	=	ND (0.33)	=	ND (0.33)	=	ND (0.33)	-	ND (0.33)
Cyclohexane	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)		ND (0.37)
1,2-Dibromo-3-chloropropane	0.02	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)
Dibromochloromethane	1	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)		ND (0.25)
1,2-Dibromoethane 1,2-Dichlorobenzene	0.03 600	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)		ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	1	ND (0.23) ND (0.16)
1,3-Dichlorobenzene 1,3-Dichlorobenzene	600	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	<del>                                     </del>	ND (0.16) ND (0.26)		ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	+	ND (0.16) ND (0.26)	-	ND (0.16) ND (0.26)	<del>                                     </del>	ND (0.16) ND (0.26)
1,4-Dichlorobenzene	75	ND (0.24)	-	ND (0.24)	-	ND (0.24)	- 1	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	- 1	ND (0.24)
Dichlorodifluoromethane	1000	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	- 1	ND (0.73)
1,1-Dichloroethane	50	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)
1,2-Dichloroethane	2	ND (0.30)	-	0.54	J -	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)	-	ND (0.30)
1,1-Dichloroethene	1	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)
cis-1,2-Dichloroethene	70	9.6	-	1.6	-	ND (0.33)	-	ND (0.33)	-	0.52	J -	ND (0.33)	-	ND (0.33)	-	ND (0.33)
trans-1,2-Dichloroethene	100	1.3	-	10.9 ND (0.43)	-	13.5 ND (0.43)	-	11.6 ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51)	-	ND (0.51)	-	ND (0.51) ND (0.43)
1,2-Dichloropropane cis-1,3-Dichloropropene	1	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)
trans-1,3-Dichloropropene	-	ND (0.32)	-	ND (0.32)	_	ND (0.32)	_	ND (0.32)	_	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)
1,4-Dioxane	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)
Ethylbenzene	700	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)
Freon 113	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)
2-Hexanone	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)
Isopropylbenzene	700	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
Methyl Acetate  Methylcyclohexane	7000	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)		ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	=	ND (3.1) ND (0.22)	<del>-</del> -	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)
Methyl Tert Butyl Ether	70	ND (0.22)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.22)	-	ND (0.26)	-	ND (0.26)	-	ND (0.22)
4-Methyl-2-pentanone(MIBK)	=	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	=	ND (1.1)	-	ND (1.1)
Methylene chloride	3	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)
Styrene 1,1,2,2-Tetrachloroethane	100	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-	ND (0.26) 2.5	-	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)	-	ND (0.26) ND (0.39)
Tetrachloroethene	1	0.43 J	J -	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	1.2	-	ND (0.35)	-	ND (0.35)	+	ND (0.35)
Toluene	600	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)
1,2,3-Trichlorobenzene	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)
1,2,4-Trichlorobenzene 1,1,1-Trichloroethane	9 30	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	+ -	ND (0.22) ND (0.32)
1,1,2-Trichloroethane	3	ND (0.32) ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.32)
Trichloroethene	1	17.5	-	0.25	J -	ND (0.25)	-	ND (0.25)	-	1.4	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)
Trichlorofluoromethane	2000	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)		ND (0.28)
Vinyl chloride m,p-Xylene	1	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	-	ND (0.17) ND (0.45)	+ -	ND (0.17) ND (0.45)
o-Xylene	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	- 1	ND (0.20)	-	ND (0.43)	-	ND (0.20)	-	ND (0.43)	- 1	ND (0.43)
Xylene (total)	1000	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)
Total VOCs	-	39.13		13.29		13.5		11.6		5.62		1.9	1	3.9	1	0
GC/MS Volatile TIC																
Total TIC, Volatile	-	0	- 1	0	- 1	0	-	0	-	0		0	- 1	0	1 - 1	0
Total Alkanes	-	0	-	0	-	0				0		0		0	-	0
Metals Analysis	1 70	076	007	1	1 1 40 1		40	1	- 40	1 40	150				1 400	
Chromium Iron	70 300	871		<10	<10 <100	<10		<10	<10 <100	<10	<10 <b>225</b>	<10	<10 <100	<10	<10 <100	-
Sodium	50000	241000		11300	++	145000		152000		16100		<10000		<10000		=
	•							•	•							
General Chemistry																
Solids, Total Dissolved Sulfate	500000 250000	1570000 527000		122000 38800	-	976000 292000	-	1050000 315000	-	151000 83300	-	<10000 <10000		<10000 <10000	1	<del>                                     </del>
	233000	52,500	1	53000		232000	1	323000	I	55500		-10000		-10000		

Sample ID		IRND4 ISCO-MW1-	IRND4 ISCO-MW1 -	IRND4 ISCO-MW1 ASC -	IRND4 ISCO-MW2	IRND4 ISCO-MW2	IRND4 ISCO-MW2 ASC	IRND4 ISCO-MW3	IRND4 ISCO-MW3	IRND4 ISCO-MW3 ASC	IRND4 ISCO-MW4	IRND4 ISCO-MW4	IRND4 ISCO-MW5	IRND4 ISCO-MW5
Lab Sample ID	NJ CLASS IIA	JB76271-15	JB76271-15F	JB76271-24	JB76271-16	JB76271-16F	JB76271-25	JB76271-17	JB76271-17F	JB76271-26	JB76271-8	JB76271-8F	JB76271-14	JB76271-14F
Sample Date	GROUNDWATER QUALITY CRITERIA (7/22/2010)	9/11/2014	9/11/2014	9/11/2014	9/11/2014	9/11/2014	9/11/2014	9/11/2014	9/11/2014	9/11/2014	9/10/2014	9/10/2014	9/11/2014	9/11/2014
Matrix	ug/L	GW	GW-FILTERED	GW	GW	GW-FILTERED	GW	GW	GW-FILTERED	GW	GW	GW-FILTERED	GW	GW-FILTERED
Unit	~ <sub>6</sub> / -	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (VOCs)					,					,		,	,	
Acetone	6000	ND (2.6)	-	ND (2.6)	304	-	235	33.8	-	27.5	ND (2.6)	-	ND (2.6)	-
Benzene	1	ND (0.21)	=	ND (0.21)	1.7	-	2.3	ND (0.21)	=	ND (0.21)	ND (0.21)	-	ND (0.21)	-
Bromochloromethane Bromodichloromethane	- 1	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	ND (0.97) ND (0.56)	-	ND (0.49) ND (0.28)	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-
Bromoform  Bromoform	1 4	ND (0.28) ND (0.31)	-	ND (0.28) ND (0.31)	ND (0.62)	-	ND (0.28) ND (0.31)	ND (0.28) ND (0.31)	-	ND (0.28) ND (0.31)	ND (0.28) ND (0.31)	-	ND (0.28) ND (0.31)	-
Bromomethane	10	ND (0.39)	-	ND (0.39)	26.7	_	32	ND (0.39)	-	ND (0.39)	ND (0.39)	_	ND (0.39)	_
2-Butanone (MEK)	300	ND (2.5)	-	ND (2.5)	239	-	185	ND (2.5)	-	ND (2.5)	ND (2.5)	-	ND (2.5)	-
Carbon disulfide	700	ND (0.50)	-	ND (0.50)	20	-	18.6	ND (0.50)	-	ND (0.50)	ND (0.50)	-	ND (0.50)	-
Carbon tetrachloride	1	ND (0.24)	-	ND (0.24)	ND (0.47)	-	ND (0.24)	ND (0.24)	-	ND (0.24)	ND (0.24)	-	ND (0.24)	-
Chlorobenzene	50	ND (0.27)	-	ND (0.27)	ND (0.54)	-	ND (0.27)	ND (0.27)	-	ND (0.27)	ND (0.27)	-	ND (0.27)	-
Chloroethane	-	ND (0.56)	-	ND (0.56)	ND (1.1)	-	ND (0.56)	ND (0.56)	-	ND (0.56)	ND (0.56)	-	ND (0.56)	-
Chloroform	70	ND (0.20)	-	ND (0.20)	3.6	-	3.3	ND (0.20)	-	ND (0.20)	0.34 J	-	ND (0.20)	-
Chloromethane	-	ND (0.33)	-	ND (0.33)	3.4	-	2.8	ND (0.33)	-	ND (0.33)	ND (0.33)	-	ND (0.33)	-
Cyclohexane	- 0.02	ND (0.37)	-	ND (0.37)	ND (0.74)	-	ND (0.37)	ND (0.37)	=	ND (0.37)	ND (0.37)	-	ND (0.37)	-
1,2-Dibromo-3-chloropropane Dibromochloromethane	0.02	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	ND (2.3) ND (0.50)	-	ND (1.2) ND (0.25)	ND (1.2) ND (0.25)	<del>                                     </del>	ND (1.2) ND (0.25)	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	<del>                                     </del>
1,2-Dibromoethane	0.03	ND (0.23)	-	ND (0.23)	ND (0.45)		ND (0.23)	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	-
1,2-Dichlorobenzene	600	ND (0.16)	-	ND (0.16)	0.68 J	-	2.7	ND (0.16)		ND (0.16)	ND (0.16)	-	ND (0.16)	-
1,3-Dichlorobenzene	600	ND (0.26)	-	ND (0.26)	ND (0.51)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
1,4-Dichlorobenzene	75	ND (0.24)		ND (0.24)	ND (0.47)	-	0.24 J	ND (0.24)	-	ND (0.24)	ND (0.24)	-	ND (0.24)	-
Dichlorodifluoromethane	1000	ND (0.73)	-	ND (0.73)	ND (1.5)	-	ND (0.73)	ND (0.73)	-	ND (0.73)	ND (0.73)	-	ND (0.73)	-
1,1-Dichloroethane	50	ND (0.35)	-	ND (0.35)	ND (0.70)	-	ND (0.35)	ND (0.35)	-	ND (0.35)	ND (0.35)	-	ND (0.35)	-
1,2-Dichloroethane	2	4	=	3.7	574	-	476	ND (0.30)	=	ND (0.30)	0.66 J	-	ND (0.30)	-
1,1-Dichloroethene	1	ND (0.50)	-	ND (0.50)	ND (0.99)	-	ND (0.50)	ND (0.50)	-	ND (0.50)	ND (0.50)	-	ND (0.50)	-
cis-1,2-Dichloroethene	70	ND (0.33)	-	ND (0.33)	ND (0.65)	-	ND (0.33)	73.8	-	66.9	ND (0.33)	-	0.50	J -
trans-1,2-Dichloroethene	100	ND (0.51)	-	ND (0.51) ND (0.43)	ND (1.0)	-	ND (0.51)	13.4 ND (0.43)	-	ND (0.51) ND (0.43)	ND (0.51)	-	ND (0.51)	-
1,2-Dichloropropane cis-1,3-Dichloropropene	1	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	ND (0.87) ND (0.57)	<u> </u>	ND (0.43) ND (0.28)	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	ND (0.43) ND (0.28)	-	ND (0.43) ND (0.28)	<del>                                     </del>
trans-1,3-Dichloropropene	-	ND (0.32)	-	ND (0.32)	ND (0.63)	_	ND (0.32)	ND (0.32)	-	ND (0.32)	ND (0.32)	_	ND (0.32)	_
1,4-Dioxane	-	ND (51)	-	ND (51)	ND (100)	-	ND (51)	ND (51)	-	ND (51)	ND (51)	-	ND (51)	-
Ethylbenzene	700	ND (0.40)	-	ND (0.40)	ND (0.79)	-	ND (0.40)	ND (0.40)	-	ND (0.40)	ND (0.40)	-	ND (0.40)	-
Freon 113	-	ND (0.45)	-	ND (0.45)	ND (0.89)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	-
2-Hexanone	-	ND (1.7)	-	ND (1.7)	ND (3.5)	-	ND (1.7)	ND (1.7)	-	ND (1.7)	ND (1.7)	-	ND (1.7)	-
Isopropylbenzene	700	ND (0.26)	=	ND (0.26)	ND (0.51)	-	ND (0.26)	ND (0.26)	=	ND (0.26)	ND (0.26)	-	ND (0.26)	-
Methyl Acetate	7000	ND (3.1)	-	ND (3.1)	ND (6.2)	-	ND (3.1)	ND (3.1)	-	ND (3.1)	ND (3.1)	-	ND (3.1)	-
Methylcyclohexane Methyl Tert Butyl Ether	70	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.43) ND (0.53)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	+
4-Methyl-2-pentanone(MIBK)	-	ND (1.1)	-	ND (1.1)	ND (0.33)	-	ND (0.20)	ND (1.1)	-	ND (0.20)	ND (1.1)	-	ND (1.1)	-
Methylene chloride	3		J -	1.4 J	2.4 J	-	2.3	ND (0.81)	-	ND (0.81)	2.1	-	ND (0.81)	-
Styrene	100	ND (0.26)	-	ND (0.26)	ND (0.51)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
1,1,2,2-Tetrachloroethane Tetrachloroethene	1	ND (0.39) ND (0.35)	-	ND (0.39) ND (0.35)	24.6 3.3	-	16.2 1.7	ND (0.39) 2.4	-	ND (0.39)	ND (0.39) <b>0.51</b> J	-	ND (0.39) <b>0.45</b>	- J -
Toluene	600	ND (0.22)	-	ND (0.33)	ND (0.44)	-	ND (0.22)	ND (0.22)	-	ND (0.22)	ND (0.22)	-	ND (0.22)	
1,2,3-Trichlorobenzene		ND (0.26)	-	ND (0.26)	ND (0.53)	-	0.68 J		-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
1,2,4-Trichlorobenzene	9	ND (0.22)	-	ND (0.22)	0.72 J	-	7.6	ND (0.22)		ND (0.22)	ND (0.22)	-	ND (0.22)	
1,1,1-Trichloroethane 1.1,2-Trichloroethane	30	ND (0.32) ND (0.28)	=	ND (0.32) ND (0.28)	ND (0.64) ND (0.55)	-	ND (0.32) 0.34 J	ND (0.32) ND (0.28)	-	ND (0.32) ND (0.28)	ND (0.32) ND (0.28)	-	0.74 ND (0.28)	J -
Trichloroethene	1	· · ·	J -	0.34 J	14.6	-	7.9	87.1	-	68.3	0.95 J		ND (0.28)	-
Trichlorofluoromethane	2000	ND (0.28)		ND (0.28)	ND (0.56)	<u>-</u>	ND (0.28)	ND (0.28)	<u>-</u>	ND (0.28)	ND (0.28)	<u>-</u>	ND (0.28)	<u> </u>
Vinyl chloride	1	ND (0.17)	-	ND (0.17)	ND (0.35)	-	ND (0.17)	ND (0.17)	-	ND (0.17)	ND (0.17)	-	ND (0.17)	-
m,p-Xylene	-	ND (0.45)	-	ND (0.45)	ND (0.90)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	-
o-Xylene Xylene (total)	1000	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	ND (0.40) ND (0.40)	-	ND (0.20) ND (0.20)	ND (0.20) ND (0.20)		ND (0.20) ND (0.20)	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	-
Total VOCs	-	6.08	1	5.44	1218.7		994.66	210.5		164.4	4.56		6.92	
														<u> </u>
GC/MS Volatile TIC														
Total TIC, Volatile	-			6.3 J			9.9 J		-		0		0	-
Total Alkanes	=	0	-	0	0	-	0	0	=	0	0	-	0	-
Metals Analysis														
Chromium	70	<10	<10	-	1280	845	-	1730	2090	- 1	55.6	<10	<10	<10
Iron	300	-		-	-	<100	-	-	<100	-	-	<del> </del>	-	835
Sodium	50000	36400	-	-	1990000	-	-	664000	-	-	13000	-	18300	-
Company Chamistan														
General Chemistry Solids, Total Dissolved	500000	245000			7120000		1	2220000	1	ı	137000	1	118000	1
Solids, Total Dissolved Sulfate	250000	192000	-	-	5280000	-	-	930000	-	-	137000 59900	-	45100	-
Ll			1	1		1			1	1	1		1	

	Sample ID NJ CLASS IIA	IRND4_ISCO-MW5 ASC	IRND4_ISCO-MW6	IRND4_ISCO-MW6	IRND4_ISCO-MW7	IRND4_ISCO-MW7	IRND4_ISCO-MW8	IRND4_ISCO-MW8	IRND4_ISCO-MW9	IRND4_ISCO-MW9	IRND4_ISCO-MW9 ASC	IRND4_IW1-DR1	IRND4_IW1-DR1	IRND4_IWI-BT2	IRND4_IWI-BT2
	Sample Date GROUNDWATER QUALITY	JB76271-23 9/11/2014	JB76271-19 9/11/2014	JB76271-19F 9/11/2014	JB76271-13 9/11/2014	JB76271-13F 9/11/2014	JB76271-20 9/11/2014	JB76271-20F 9/11/2014	JB76271-10 9/10/2014	JB76271-10F 9/10/2014	JB76271-22 9/10/2014	JB76271-7 9/10/2014	JB76271-7F 9/10/2014	JB76271-18 9/11/2014	JB76271-18F 9/11/2014
	Matrix	GW	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED	5/15/2014 GW	GW	GW-FILTERED	GW	GW-FILTERED
	ug/L Unit	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Volatile Organic Compounds (V	VOCs)														
Acetone	6000	ND (2.6)	ND (2.6)	=	45	-	47.4	-	ND (2.6)	-	9.2 J	112 (210)	=	9.5	J -
Benzene Bromoshloromothana	1	ND (0.21) ND (0.49)	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	-	ND (0.21) ND (0.49)	ND (0.21)	-	ND (0.21)	-
Bromochloromethane Bromodichloromethane	1	ND (0.49) ND (0.28)	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	ND (0.49) ND (0.28)	-	ND (0.49) ND (0.28)	-
Bromoform	4	ND (0.31)	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	-	ND (0.31)	ND (0.31)	-	ND (0.31)	-
Bromomethane	10	ND (0.39)	ND (0.39)	-	13.7	-	ND (0.39)	-	ND (0.39)	-	ND (0.39)	ND (0.39)	-	ND (0.39)	-
2-Butanone (MEK)	300	ND (2.5)	ND (2.5)	=	ND (2.5)	-	ND (2.5)	-	ND (2.5)	=	ND (2.5)	ND (2.5)	=	ND (2.5)	-
Carbon disulfide Carbon tetrachloride	700 1	ND (0.50) ND (0.24)	ND (0.50) ND (0.24)	-	ND (0.50) <b>0.73</b> J	-	ND (0.50) <b>0.31</b>	-	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	ND (0.50) ND (0.24)	-	ND (0.50) ND (0.24)	-
Chlorobenzene	50	ND (0.24)	ND (0.24)	-	ND (0.27)	-	ND (0.27)	-	ND (0.24) ND (0.27)	-	ND (0.24) ND (0.27)	ND (0.24)	-	ND (0.24)	-
Chloroethane	-	ND (0.56)	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	-	ND (0.56)	ND (0.56)	-	ND (0.56)	-
Chloroform	70	ND (0.20)	ND (0.20)	Ξ	1.1	ē	1.1	=	ND (0.20)	=	ND (0.20)	ND (0.20)	=	ND (0.20)	=
Chloromethane	-	ND (0.33)	ND (0.33)	-	2.8	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	-	ND (0.33)	-
Cyclohexane	0.02	ND (0.37)	ND (0.37) ND (1.2)	-	ND (0.37) ND (1.2)	-	ND (0.37)	-	ND (0.37) ND (1.2)	-	ND (0.37)	ND (0.37)	-	ND (0.37) ND (1.2)	+ -
1,2-Dibromo-3-chloropropane Dibromochloromethane	0.02	ND (1.2) ND (0.25)	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	ND (1.2) ND (0.25)	-	ND (1.2) ND (0.25)	+
1,2-Dibromoethane	0.03	ND (0.23)	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	-	ND (0.23)	ND (0.23)	-	ND (0.23)	-
1,2-Dichlorobenzene	600	ND (0.16)	ND (0.16)	-	ND (0.16)		ND (0.16)	-	ND (0.16)	-	ND (0.16)	ND (0.16)	-	ND (0.16)	-
1,3-Dichlorobenzene	600	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
1,4-Dichlorobenzene Dichlorodifluoromethane	75 1000	ND (0.24) ND (0.73)	ND (0.24) ND (0.73)	-	ND (0.24) ND (0.73)	-	ND (0.24) ND (0.73)	-	ND (0.24) ND (0.73)	-	ND (0.24) ND (0.73)	ND (0.24) ND (0.73)	-	ND (0.24) ND (0.73)	-
1,1-Dichloroethane	50	ND (0.75)	ND (0.75) ND (0.35)	-	ND (0.75)	-	ND (0.75) ND (0.35)	-	ND (0.75) ND (0.35)	-	ND (0.75) ND (0.35)	ND (0.75) ND (0.35)	-	ND (0.75)	<del>                                     </del>
1,2-Dichloroethane	2	ND (0.30)	2.4	-	4.6	-	1.7	-	ND (0.30)	-	0.79 J	ND (0.30)	-	ND (0.30)	-
1,1-Dichloroethene	1	ND (0.50)	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	ND (0.50)	-	ND (0.50)	-
cis-1,2-Dichloroethene	70	0.84	J 3	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	ND (0.33)	-	4.8	-
trans-1,2-Dichloroethene 1,2-Dichloropropane	100	ND (0.51) ND (0.43)	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	-	ND (0.51) ND (0.43)	=	ND (0.51) ND (0.43)	ND (0.51) ND (0.43)	-	0.55 ND (0.43)	-
cis-1,3-Dichloropropene	-	ND (0.43) ND (0.28)	ND (0.43) ND (0.28)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43) ND (0.28)	ND (0.43) ND (0.28)	-	ND (0.43)	<del>                                     </del>
trans-1,3-Dichloropropene	-	ND (0.32)	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	ND (0.32)	-	ND (0.32)	-
1,4-Dioxane	-	ND (51)	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	-	ND (51)	ND (51)	-	ND (51)	-
Ethylbenzene	700	ND (0.40)	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	-	ND (0.40)	ND (0.40)	-	ND (0.40)	-
Freon 113 2-Hexanone		ND (0.45) ND (1.7)	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)	ND (0.45) ND (1.7)	-	ND (0.45) ND (1.7)	-
Isopropylbenzene	700	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
Methyl Acetate	7000	ND (3.1)	ND (3.1)	=	ND (3.1)	÷	ND (3.1)	-	ND (3.1)	=	ND (3.1)	ND (3.1)	=	ND (3.1)	-
Methylcyclohexane	-	ND (0.22)	ND (0.22)	=	ND (0.22)	-	ND (0.22)	-	ND (0.22)	=	ND (0.22)	ND (0.22)	=	ND (0.22)	-
Methyl Tert Butyl Ether 4-Methyl-2-pentanone(MIBK)	70	ND (0.26) ND (1.1)	ND (0.26) ND (1.1)	-	ND (0.26) ND (1.1)	-	ND (0.26) ND (1.1)	-	ND (0.26) ND (1.1)	-	ND (0.26) ND (1.1)	ND (0.26) ND (1.1)	-	ND (0.26) ND (1.1)	-
Methylene chloride	3	ND (0.81)	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	-	ND (0.81)	ND (0.81)	-	ND (0.81)	-
Styrene	100	ND (0.26)	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	ND (0.26)	-	ND (0.26)	-
1,1,2,2-Tetrachloroethane Tetrachloroethene	1 1	ND (0.39) <b>0.41</b>	ND (0.39) J ND (0.35)	-	ND (0.39) <b>0.63</b> J	-	ND (0.39) <b>0.42</b>	J -	ND (0.39) 1.4	-	ND (0.39) 1	ND (0.39) <b>0.57</b> J	-	ND (0.39) 0.42	· ·
Toluene	600	ND (0.22)	ND (0.22)	=	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	ND (0.22)	=	ND (0.22)	-
1,2,3-Trichlorobenzene 1,2,4-Trichlorobenzene	- 9	ND (0.26) ND (0.22)	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)		ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	-	ND (0.26) <b>0.47</b> J	ND (0.26) ND (0.22)	-	ND (0.26) ND (0.22)	-
1,1,1-Trichloroethane	30	0.85	J ND (0.22)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)		ND (0.32)	ND (0.22) ND (0.32)	-	ND (0.22) ND (0.32)	
1,1,2-Trichloroethane	3	ND (0.28)	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	ND (0.28)	-	ND (0.28)	-
Trichloroethene Trichlorofluoromethane	2000	4.4 ND (0.28)	5.4 ND (0.28)	-	1.4 ND (0.28)	-	1.2 ND (0.28)	-	14.8 ND (0.28)	-	13.3 ND (0.28)	2.4 ND (0.28)	-	12.7 ND (0.28)	-
Vinyl chloride	1	ND (0.17)	ND (0.28) ND (0.17)	-	ND (0.28) ND (0.17)	-	ND (0.28) ND (0.17)	-	ND (0.28) ND (0.17)	-	ND (0.28) ND (0.17)	ND (0.28) ND (0.17)	-	ND (0.28) ND (0.17)	-
m,p-Xylene	-	ND (0.45)	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	-	ND (0.45)	ND (0.45)	-	ND (0.45)	-
o-Xylene Xylene (total)	1000	ND (0.20) ND (0.20)	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	ND (0.20) ND (0.20)	-	ND (0.20) ND (0.20)	+ -
Total VOCs	-	6.5	10.8		69.96				16.2		24.76	2.97		27.97	
CO (NACA VIII A TIC															
GC/MS Volatile TIC Total TIC, Volatile		0		_ 1	24 J	1 -	1 0 1		0		0	0		0	
Total Alkanes		0	0	-	0	-	0	-	0	-	0	0	-	0	-
				<u> </u>					<u> </u>						
Metals Analysis	70			<10	816	007	3210	2110	10		1		1 410	611	1 <10
Chromium Iron	70 300	-	<10	<10 <100	816	907 227		2110 1360	10 -	<10 <100	-	<10		611	<10 <100
Sodium	50000	-	92300	-	1530000	-	1890000	-	42500	-	-	14500		232000	-
General Chemistry															
Solids, Total Dissolved	500000	-	394000	- 1	6860000	- 1	5930000		460000	- 1	- 1	96700	1	1160000	- 1
	250000		211000		2720000		1140000	+	124000	+	t	96800		444000	+

Sample Date	N) CLASS IIA JNDWATER QUALITY TERIA (7/22/2010) ug/L  6000 1 - 1 4 10 300 700	IRND4_MW-10S JB76271-5 9/10/2014 GW ug/L  ND (2.6) ND (0.21) ND (0.49) ND (0.28) ND (0.31)	IRND4_MW-10S JB76271-5F 9/10/2014 GW-FILTERED ug/L -	IRND4_DUP JB76271-6 9/10/2014 GW ug/L ND (2.6) ND (0.21)	IRND4_DUP JB76271-6F 9/10/2014 GW-FILTERED ug/L	IRND4_MW-14SD JB76271-3 9/10/2014 GW ug/L	IRND4_MW-14SD JB76271-3F 9/10/2014 GW-FILTERED ug/L	IRND4_MW-14SS JB76271-2 9/10/2014 GW ug/L	IRND4_MW-14SS JB76271-2F 9/10/2014 GW-FILTERED	IRND4_MW-5I JB76271-4 9/10/2014 GW	IRND4_MW-5I JB76271-4F 9/10/2014 GW-FILTERED	IRND4_MW11I JB76271-1 9/10/2014 GW	IRND4_MW11I JB76271-1F 9/10/2014 GW-FILTERED	IRND4_PZ-1S JB76271-9 9/10/2014 GW	IRND4_P2-1S JB76271-9F 9/10/2014 GW-FILTERED
Sample Date Matrix Unit Volatile Organic Compounds (VOCs)  Acetone Benzene Bromochloromethane Bromodichloromethane Bromoform Bromomethane	6000 1 - 1 - 1 1 1 1 1 1 1 1 1 1 1 1 1 1	MD (2.6) ND (0.21) ND (0.49) ND (0.28) ND (0.31)	GW-FILTERED ug/L	GW ug/L ND (2.6) ND (0.21)	GW-FILTERED ug/L	GW	GW-FILTERED	GW	GW-FILTERED						
Matrix Unit Volatile Organic Compounds (VOCs)  Acetone Benzene Bromochloromethane Bromodichloromethane Bromoform Bromomethane	1 1 4 10 300 700	ND (2.6) ND (0.21) ND (0.49) ND (0.28) ND (0.31)	ug/L	ug/L  ND (2.6)  ND (0.21)	ug/L					GW	GW-FILTERED	GW	GW-FILTERED	GW	GW-FILTERED
Volatile Organic Compounds (VOCs)  Acetone Benzene Bromochloromethane Bromodichloromethane Bromoform Bromomethane	1 - 1 4 10 300 700	ND (2.6) ND (0.21) ND (0.49) ND (0.28) ND (0.31)	,	ND (2.6) ND (0.21)	,	ug/L	ug/L								
Acetone Benzene Bromochloromethane Bromodichloromethane Bromoform Bromomethane	1 - 1 4 10 300 700	ND (0.21) ND (0.49) ND (0.28) ND (0.31)		ND (0.21)	- 1			ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Benzene Bromochloromethane Bromodichloromethane Bromoform Bromomethane	1 - 1 4 10 300 700	ND (0.21) ND (0.49) ND (0.28) ND (0.31)	-	ND (0.21)		ND (2.6)		ND (2.6)		ND (2.6)		ND (2.6)	1 . 1	ND (2.6)	I - I
Bromodichloromethane Bromoform Bromomethane	4 10 300 700	ND (0.28) ND (0.31)	-		-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-	ND (0.21)	-
Bromoform Bromomethane	4 10 300 700	ND (0.31)	_	ND (0.49)	=	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-	ND (0.49)	-
Bromomethane	10 300 700			ND (0.28)	=	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-
	300 700		-	ND (0.31)	-	ND (0.31) ND (0.39)	-	ND (0.31)	-	ND (0.31) ND (0.39)	-	ND (0.31) ND (0.39)	-	ND (0.31) ND (0.39)	-
	700	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	-	ND (0.39) ND (2.5)	+ -
Carbon disulfide		ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-	ND (0.50)	-
Carbon tetrachloride	1	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-
Chlorobenzene	50	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-	ND (0.27)	-
Chloroethane Chloroform	70	ND (0.56) ND (0.20)	-	ND (0.56) ND (0.20)	-	ND (0.56) 0.29	-	ND (0.56) <b>0.27</b> J	-	ND (0.56) ND (0.20)	-	ND (0.56) 0.21		ND (0.56) ND (0.20)	<del>-</del> -
Chloromethane	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	_	ND (0.33)	-	ND (0.33)	-
Cyclohexane	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-	ND (0.37)	-
1,2-Dibromo-3-chloropropane	0.02	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)	-	ND (1.2)		ND (1.2)		ND (1.2)	
Dibromochloromethane	1	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	-	ND (0.25)	
1,2-Dibromoethane 1,2-Dichlorobenzene	0.03 600	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	1 -	ND (0.23) ND (0.16)	-	ND (0.23) ND (0.16)	<del>                                     </del>	ND (0.23) ND (0.16)		ND (0.23) ND (0.16)	+ -
1,3-Dichlorobenzene	600	ND (0.16)	-	ND (0.26)	-	ND (0.26)	-	ND (0.16)	-	ND (0.16)	-	ND (0.26)	-	ND (0.26)	+ - +
1,4-Dichlorobenzene	75	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	-	ND (0.24)	- 1
Dichlorodifluoromethane	1000	ND (0.73)	-	ND (0.73)	-	ND (0.73)		ND (0.73)	-	ND (0.73)	-	ND (0.73)	-	ND (0.73)	
1,1-Dichloroethane	50	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	<del>                                     </del>
1,2-Dichloroethane 1,1-Dichloroethene	1	<b>3.1</b> ND (0.50)	-	3 ND (0.50)	-	ND (0.30) ND (0.50)	1 1	ND (0.30) ND (0.50)	-	1.3 ND (0.50)	-	ND (0.30) ND (0.50)	-	5.3 ND (0.50)	+ - +
cis-1,2-Dichloroethene	70	0.76	J -	0.92 J	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	-	ND (0.33)	+ - +
trans-1,2-Dichloroethene	100	ND (0.51)		ND (0.51)	<u>-</u>	ND (0.51)	<u> </u>	ND (0.51)		ND (0.51)		ND (0.51)		ND (0.51)	-
1,2-Dichloropropane	1	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	-	ND (0.43)	
cis-1,3-Dichloropropene	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	<del>                                     </del>
trans-1,3-Dichloropropene 1,4-Dioxane	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	-	ND (0.32) ND (51)	+ -
Ethylbenzene	700	ND (51) ND (0.40)	-	ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	-	ND (51) ND (0.40)	+ -
Freon 113		ND (0.45)		ND (0.45)	<u>-</u>	ND (0.45)	<u> </u>	ND (0.45)		ND (0.45)		ND (0.45)		ND (0.45)	<u> </u>
2-Hexanone	-	ND (1.7)	-	ND (1.7)	-	ND (1.7)	1	ND (1.7)	-	ND (1.7)		ND (1.7)	-	ND (1.7)	
Isopropylbenzene Mothyl Acotato	700	ND (0.26)	-	ND (0.26)	-	ND (0.26)	1 -	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	+
Methyl Acetate Methylcyclohexane	7000	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-	ND (3.1) ND (0.22)	-
Methyl Tert Butyl Ether	70	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	-	ND (0.26)	
4-Methyl-2-pentanone(MIBK)	-	ND (1.1)	-	ND (1.1)	=	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-	ND (1.1)	-
Methylene chloride Styrene	3 100	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	-	ND (0.81) ND (0.26)	<del>-</del>
1,1,2,2-Tetrachloroethane	1	ND (0.39)		ND (0.39)	<u>-</u>	ND (0.39)	<u> </u>	ND (0.39)		ND (0.39)		ND (0.39)		ND (0.39)	
Tetrachloroethene	1	1	-		-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	-	ND (0.35)	
Toluene 1,2,3-Trichlorobenzene	- 600	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)	-	ND (0.22) ND (0.26)		ND (0.22) ND (0.26)	+ - +
1,2,4-Trichlorobenzene	9	ND (0.22)	-	ND (0.22)	-	ND (0.22)	- 1	ND (0.22)	-	ND (0.22)	-	ND (0.22)	-	ND (0.22)	
1,1,1-Trichloroethane	30	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	ND (0.32)	-	0.6	-	ND (0.32)	
1,1,2-Trichloroethane Trichloroethene	3 1	ND (0.28) 1.5	-	ND (0.28) 1.5	-	ND (0.28) ND (0.25)	-	ND (0.28) ND (0.25)	-	ND (0.28) <b>0.27</b> J	-	ND (0.28) ND (0.25)	-	ND (0.28)	+ - +
Trichlorofluoromethane	2000	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	-	ND (0.28)	
Vinyl chloride	1	ND (0.17)	-	ND (0.17)	-	ND (0.17)	<u> </u>	ND (0.17)	-	ND (0.17)		ND (0.17)		ND (0.17)	<del>                                     </del>
m,p-Xylene o-Xylene	-	ND (0.45) <b>0.27</b>	1 - -	ND (0.45) 0.31 J	-	ND (0.45) ND (0.20)	<del>                                     </del>	ND (0.45) ND (0.20)		ND (0.45) ND (0.20)		ND (0.45) ND (0.20)		ND (0.45) ND (0.20)	+
Xylene (total)	1000	0.51	J -	0.53 J		ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	-	ND (0.20)	
Total VOCs	-	6.87		6.86		0.29	1	0.27		1.57		0.81		7.3	
GC/MS Volatile TIC															
Total TIC, Volatile	-	0	-	0	-	0	-	0	-		-	0	-	0	<u> </u>
Total Alkanes	-	0	-	0	-	0	-	0	-	0	-	0	-	0	-
Metals Analysis															
Chromium	70	24.7	12.2	23.6	11.7	<10	<10	<10	<10	<10	<10	<10	<10	31.6	17.9
Iron	300	-	<100	-	<100	-	<100	-	<100	-	<100	-	<100		146
Sodium	50000	43000	-	37800	-	58000	-	61000	-	<10000	-	<10000	-	71300	-
General Chemistry															
Solids, Total Dissolved	500000	383000		400000	-	376000		416000	-	50000	-	62000	-	475000	-
Sulfate	250000	171000	=	176000	-	101000	-	117000	-	68300	-	63800	-	101000	=

# Attachment 3 Data Validation Results

TO: J. Levesque cc:

**FROM:** K. Storne

**RE:** Evor Phillips Leasing Company (EPLC)Superfund Site, OU3

Site Groundwater RA Baseline Sampling Event, Data

Validation Report

**FILE:** 19726/51308.015.100

**DATE:** June 04, 2015

This report presents the data validation results performed for environmental samples collected in March 2015 for the 2015 Annual/Baseline Groundwater Monitoring Event as part of the OU3-Site Groundwater Remedial Action at the Evor Phillips Leasing Company (EPLC) Superfund Site in Old Bridge Township, New Jersey.

### **SAMPLE AND VALIDATION SUMMARY**

The environmental samples collected for this effort consisted of groundwater samples, matrix spike/matrix spike duplicates, field duplicate, field blanks and trip blanks. Samples were analyzed by Accutest Laboratories of Dayton, New Jersey (Accutest New Jersey).

The laboratory utilized the methods listed in Table 1 for sample analyses.

Table 1. Analytical methods and	references	
Parameter	Methods	Reference
VOCs	USEPA Methods8000C/5030B/8260B	1
Metals	USEPA Methods 3010A/6010C	2
Sulfate	USEPA Method 9056A/300.0	2/4
TDS	SM20 2540C	3

### Note:

VOCs indicates volatile organic compounds.

TDS indicates total dissolved solids.

- 1. United States Environmental Protection Agency (USEPA). 2004. *Test Methods for Evaluating Solid Waste: Physical/Chemical Methods, SW-846,* 3rd Edition, Update IIIB. Washington D.C.
- 2. USEPA. 2007. Test Methods for Evaluating Solid Waste: Physical/Chemical Methods, SW-846, 3rd Edition, Update IV. Washington D.C.
- 3. AWWA, APHA, WEF. 1998. Standard Methods for the Examination of Water and Wastewater, 20th Edition. Washington, D.C.
- 4. USEPA. 1993a. Methods for the Determination of Inorganic Substances in Environmental Samples, EPA-600/R-93/100. Washington, D.C.

The laboratory data packages included summary forms for quality control analysis and supportive raw data.

The samples submitted for data review are summarized in the attached Table 2. Table 3 presents the specific data validation approach applied to data generated. Table 4 presents the Laboratory quality assurance/quality control (QA/QC) analyses definitions.

In accordance with the approved RDR/RAWP, full validation was performed on 10 percent of the samples collected and submitted for validation. This consisted of a review of data summary forms and raw analytical data provided in the data packages. Partial validation was performed for the remaining data. Partial data quality review consists of a review of only analytical QC summary forms that are included in the data package. The forms and the information contained on the forms are not evaluated for accuracy or completeness during partial data validation.

The analytical data generated for this investigation were evaluated by O'Brien & Gere using the QA/QC criteria established in the methods utilized by the laboratories and the following document:

 O'Brien & Gere. 2014. Uniform Federal Policy Quality Assurance Project Plan, Operable Unit 3 (OU3)- Site Groundwater Evor Phillips Leasing Company (EPLC) Superfund Site, Old Bridge Township, New Jersey. Edison, New Jersey. (QAPP)



Data affected by excursions from these criteria were qualified using professional judgment and the general validation approach provided in the following validation guideline documents, modified to reflect the requirements of the methods utilized by the laboratories:

- New Jersey Department of Environmental Protection (NJDEP). 2001a. Standard Operating Procedure (SOP) for Analytical Data Validation of Target Analyte List (TAL) – Inorganics, SOP No. 5.A.2. Trenton, New Jersey
- NJDEP. 2001b. Standard Operating Procedures for the Quality Assurance Data Validation of Analytical Deliverables TCL- Organics (based on the USEPA SOW OLM04.2 with Revisions), SOP No. 5.A.13. Trenton, New Jersey

The application of these validation guidelines has been modified to reflect the requirements of the methods utilized by the laboratory.

In accordance with the NJDEP guidance, and utilizing professional judgment, the following qualifiers are used in this type of data review:

- "U" Indicates that the analyte was analyzed for, but was not detected.
- "J" Indicates that the result should be considered to be an estimated value. This qualifier is used when the data validation process identifies a deficiency in the data generation process.
- "UJ" Indicates that the sample-specific reporting limit for the analyte in this sample should be considered approximate. This qualifier is used when the data validation process identifies a deficiency in the data generation process.
- "R" Indicates that the reporting limit or sample result has been determined to be unusable due to a major deficiency in the data generation process. The data should not be used for any qualitative or quantitative purposes.

In addition, in accordance with the NJDEP guidance, the following single word descriptors were added to analyte results if the reported analyte required a quality assurance action.

- Qualify (Q) used when the results of a given analyte in a sample do not meet all QA/QC criteria but the deficiencies are not severe enough to warrant data rejection.
- Negate (N) used when the presence of a given analyte in a sample can be attributed to the laboratory/field introduced contamination.
- Reject (R) used when the results of a given analyte in a sample do not meet all QA/QC criteria so that the
  qualitative presence and/or quantitation of that analyte in the sample cannot be determined with any
  degree of confidence.

Footnotes, based on the NJDEP validation guidance, were applied to each qualifier to define the type of excursion that affected the sample result, resulting in the qualification of the data. The footnote used in this validation is presented in Table 5 below.



Table 5. Validation Footnote Definitions	
Footnote	Type of Excursion
39	The reported concentration is quantitative qualified because the concentration is below the RL.

The following parameters were evaluated, where applicable:

- QAPP compliance
- Documentation completeness
- · Chain-of-custody record
- Sample collection
- Sample preservation
- Holding times
- Calibrations (Full validation only)
- Blank analysis
- Matrix spike/ matrix spike duplicate (MS/MSD) analysis
- Laboratory Control Sample (LCS) analysis
- Field duplicate analysis
- Surrogate recovery
- Internal standards performance
- Gas chromatography/mass spectrometry (GC/MS) instrument performance check (Full validation only)
- Inductively coupled plasma (ICP) interference check analysis (Full validation only)
- ICP serial dilution analysis
- Laboratory duplicate analysis
- Target analyte quantitation, identification, and quantitation limits (QLs) (Full validation only)

The following sections of this memorandum present the results of the comparison of the analytical data to the QA/QC criteria specified above.

### CHAIN OF CUSTODY RECORDS AND SAMPLE COLLECTION

For samples collected 3/4/15 to 3/10/15, the sample collection dates were not documented for the complete list of samples on the record. However, the collection date was included in the sample identification.

Although the record for 2BASE\_ISCO-MW5\_03092015 only listed VOC analysis, sample containers were also submitted for metal, TDS, and sulfate analyses. This revision was confirmed by O'Brien & Gere on 3/12/15.

A field duplicate sample was not submitted for total and dissolved metal, TDS and sulfate analyses.

### **VOC DATA EVALUATION SUMMARY**

The following QA/QC parameters were found to meet method and validation criteria or did not result in additional qualification of sample results:

- QAPP compliance
- Documentation completeness
- Sample preservation



- Holding times
- Calibrations (Full validation only)
- Blank analysis
- MS/MSD analysis
- LCS analysis
- Field duplicate analysis
- Surrogate recovery
- Internal standards performance
- GC/MS instrument performance check (Full validation only)
- Target analyte identification

Excursions from method or validation criteria were not identified during the validation process. Additional observations are described below.

### I. Target analyte quantitation and detection limits

Sample results with concentrations greater than the method detection limits (MDL) but less than the QL were flagged as approximate (J) by the laboratory. This flag was retained during the validation process to indicate the data is approximate (J, 39).

### METALS, SULFATE and TDS DATA EVALUATION SUMMARY

The following QA/QC parameters were found to meet method and validation criteria or did not result in additional qualification of sample results (where applicable):

- QAPP compliance
- Documentation completeness
- Sample preservation
- Holding times
- Calibrations (Full validation only)
- Blank analysis
- MS/MSD analysis
- LCS analysis
- ICP interference check analysis (Full validation only)
- ICP serial dilution analysis
- Laboratory duplicate analysis

Excursions from method or validation criteria were not identified during the validation process. Additional observations are described below.

### I. Field duplicate analysis

A field duplicate sample was not submitted for total and dissolved metal, TDS and sulfate analyses. Therefore, field precision could not be evaluated for these analyses.

### II. Target analyte quantitation and QLs

Results for metals and inorganics were reported to the QL concentration.



Metals and sulfate were reported using dilutions due to elevated concentrations of target analytes.

### **DATA USABILITY**

The data from the samples on Table 2 were evaluated based on QA/QC criteria established by the methods listed in Table 1 and the data validation approach as described in Table 3.

Major deficiencies in the data generation process would have resulted in data points being rejected, indicating that the data are considered unusable for either quantitative or qualitative purposes. Major deficiencies were not identified during the validation process. Minor deficiencies in the data generation process would have resulted in sample data being characterized as approximate or non-detected. Minor deficiencies were not identified during the validation process.

A discussion of the data quality with regard to the data usability parameters follows:

<u>Precision</u>: Data were not rejected for precision excursions.

<u>Sensitivity</u>: Sensitivity is established by QLs, which represent measurable concentrations of analytes which can be determined with a designated level of confidence, that meet project requirements. Dilutions were performed for analyses due to elevated concentrations of target analytes in the samples.

<u>Accuracy</u>: Data were not rejected for accuracy excursions.

Representativeness: Data were not rejected for representativeness excursions.

<u>Comparability</u>: Data usability with respect to comparability is 100 percent, as standardized analytical methods, QLs, reference materials, and data deliverables were used throughout the data generation process for this project.

<u>Completeness</u>: For the samples submitted for data validation, overall data usability with respect to completeness 100 percent for the data, considering the complete data set; therefore, the usability met the QAPP requirement of usable for qualitative and quantitative purposes.



Table 2. Sample Cross Reference Table

Samples collected and submitted for data validation

I ahawatawa Nama	Data Callagted	Client Identification	Laboratory	Matrix	Analysis Resussated
Laboratory Name		Client Identification	Identification		Analysis Requested
Accutest	3/4/2015	2BASE_WCC-3M_03042015	JB89329-1	Groundwater	VOCs
Accutest	3/4/2015	2BASE_WCC-1M_03042015	JB89329-2	Groundwater	VOCs
Accutest	3/4/2015	2BASE_WCC-1S_03042015	JB89329-3	Groundwater	VOCs
Accutest	3/4/2015	2BASE_MW-15D_03042015	JB89329-4	Groundwater	VOCs
Accutest	3/4/2015	2BASE_DUP_03042015[2BASE_WCC-3M_03042015]	JB89329-5	Groundwater	VOCs
Accutest	3/4/2015	2BASE_MW-19S_03042015	JB89329-6	Groundwater	VOCs
Accutest	3/4/2015	2BASE_MW-9I_03042015	JB89329-7	Groundwater	VOCs
Accutest	3/4/2015	2BASE_MW-11I_03042015	JB89329-8	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/4/2015	2BASE_MW-11I_03042015-FILTERED	JB89329-8F	Groundwater	Dissolved Metals
Accutest	3/4/2015	2BASE_MW-5I_03042015	JB89329-9	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/4/2015	2BASE_MW-5I_03042015-FILTERED	JB89329-9F	Groundwater	Dissolved Metals
Accutest	3/4/2015	2BASE_IW1-DR1_03042015	JB89329-10	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/4/2015	2BASE_IW1-DR1_03042015-FILTERED	JB89329-10F	Groundwater	Dissolved Metals
Accutest	3/4/2015	2BASE_PZ-1S_03042015	JB89329-11	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/4/2015	2BASE_PZ-1S_03042015-FILTERED	JB89329-11F	Groundwater	Dissolved Metals
Accutest	3/4/2015	2BASE_MW-24_03042015	JB89329-12	Groundwater	VOCs
Accutest	3/4/2015	2BASE_TB_03042015	JB89329-13	Aqueous	VOCs
Accutest	3/6/2015	2BASE_ISCO-MW3_03062015	JB89329-14	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE_ISCO-MW3_03062015-FILTERED	JB89329-14F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE IW1-BT2 03062015	JB89329-15	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE IW1-BT2 03062015-FILTERED	JB89329-15F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE ISCO-MW8 03062015	JB89329-16	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE ISCO-MW8 03062015-FILTERED	JB89329-16F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE IW-4S 03062015, MS/MSD	JB89329-17	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE IW-4S 03062015-FILTERED, MS/MSD	JB89329-17F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE ISCO-MW6 03062015	JB89329-18	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE ISCO-MW6 03062015-FILTERED	JB89329-18F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE ISCO-MW7 03062015	JB89329-19	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE ISCO-MW7_03062015-FILTERED	JB89329-19F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE ISCO-MW2 03062015	JB89329-20	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE ISCO-MW2 03062015-FILTERED	JB89329-20F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE ISCO-MW4 03062015	JB89329-201 JB89329-21	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE ISCO-MW4_03062015-FILTERED	JB89329-21F	Groundwater	Dissolved Metals
Accutest	3/6/2015	2BASE FB 03062015	JB89329-21	Aqueous	VOCs, Metals, Sulfate, TDS
Accutest	3/6/2015	2BASE FB 03062015-FILTERED	JB89329-22F	Aqueous	Dissolved Metals
Accutest	3/6/2015	2BASE_FB_03062015 2BASE_TB_03062015	JB89329-23	Aqueous	VOCs
Accutest	3/9/2015	2BASE_ISCO-MW1_03092015		Groundwater	VOCs, Metals, Sulfate, TDS
	3/9/2015		JB89329-24	Groundwater	
Accutest		2BASE_ISCO-MW1_03092015-FILTERED	JB89329-24F		Dissolved Metals VOCs
Accutest	3/9/2015	2BASE_MW-6S_03092015 2BASE_ISCO-MW5_03092015	JB89329-25 JB89329-26	Groundwater	
Accutest	3/9/2015			Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/9/2015	2BASE_ISCO-MW5_03092015-FILTERED	JB89329-26F	Groundwater	Dissolved Metals
Accutest	3/9/2015	2BASE_MW-23D_03092015	JB89329-27	Groundwater	VOCs
Accutest	3/9/2015	2BASE_MW-23S_03092015	JB89329-28	Groundwater	VOCs
Accutest	3/9/2015	2BASE_MW-23I_03092015	JB89329-29	Groundwater	VOCs
Accutest	3/9/2015	2BASE_ISCO-MW9_03092015	JB89329-30	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/9/2015	2BASE_ISCO-MW9_03092015-FILTERED	JB89329-30F	Groundwater	Dissolved Metals
Accutest	3/9/2015	2BASE_MW-28_03092015	JB89329-31	Groundwater	VOCs
Accutest	3/9/2015	2BASE_FB_03092015	JB89329-32	Aqueous	VOCs, Metals, Sulfate, TDS
Accutest	3/9/2015	2BASE_FB_03092015-FILTERED	JB89329-32F	Aqueous	Dissolved Metals
Accutest	3/9/2015	2BASE_TB_03092015	JB89329-33	Aqueous	VOCs
Accutest	3/10/2015	2BASE_EW-3_03102015, MS/MSD	JB89329-34	Groundwater	VOCs
Accutest	3/10/2015	2BASE_MW-14SS_03102015	JB89329-35	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/10/2015	2BASE_MW-14SS_03102015-FILTERED	JB89329-35F	Groundwater	Dissolved Metals
Accutest	3/10/2015	2BASE_MW-14SD_03102015	JB89329-36	Groundwater	VOCs, Metals, Sulfate, TDS
Accutest	3/10/2015	2BASE_MW-14SD_03102015-FILTERED	JB89329-36F	Groundwater	Dissolved Metals
Accutest	3/10/2015	2BASE_MW-10S_03102015	JB89329-37	Groundwater	VOCs, Metals, Sulfate, TDS
			JB89329-37F	Groundwater	Dissolved Metals

Note

Accutest indicates Accutest Laboratories of Dayton, New Jersey.

VOCs indicates volatile organic compounds.

TDS indicates total dissolved solids.

MS/MSD indicates matrix spike/matrix spike duplicate.

DUP indicates field duplicate.

The sample identification utilized for field duplicate is shown in brackets.

TB indicates trip blank.

FB indicates field blank.

### Table 3 - O'Brien & Gere data validation approach using NJDEP data validation guidelines

### General Validation Approach

Data evaluation is based on QA/QC criteria established the methods utilized by the laboratory and quality plans developed for the project.

The NJDEP data validation guidance applies to data generated using USEPA CLP methods. This project was not analyzed using CLP methods. Therefore, data affected by excursions from criteria presented in the methods and quality plan are qualified using professional judgment with some consideration of the general guidance provided in the following documents:

- New Jersey Department of Environmental Protection (NJDEP). 2001a. Standard Operating Procedures for the Quality
  Assurance Data Validation of Analytical Deliverables TCL- Organics (based on the USEPA SOW OLM04.2 with Revisions),
  SOP No. 5.A.13. Trenton, New Jersey; and
- NJDEP. 2001b. Standard Operating Procedure (SOP) for Analytical Data Validation of Target Analyte List (TAL) Inorganics, SOP No. 5.A.2. Trenton, New Jersey.

The following qualifiers are applied to data:

""U" Indicates that the analyte was analyzed for, but was not detected.

"J" Indicates that the result should be considered to be an estimated value. This qualifier is used when the data validation process identifies a deficiency in the data generation process.

"UJ" Indicates that the sample-specific reporting limit for the analyte in this sample should be considered approximate. This qualifier is used when the data validation process identifies a deficiency in the data generation process.

"R" Indicates that the reporting limit or sample result has been determined to be unusable due to a major deficiency in the data generation process. The data should not be used for any qualitative or quantitative purposes.

In addition, in accordance with the NJDEP guidance, the following single word descriptors were added to analyte results if the reported analyte required a quality assurance action.

- Qualify (Q) used when the results of a given analyte in a sample do not meet all QA/QC criteria but the deficiencies are not severe enough to warrant data rejection.
- Negate (N) used when the presence of a given analyte in a sample can be attributed to the laboratory/field introduced contamination.
- Reject (R) used when the results of a given analyte in a sample do not meet all QA/QC criteria so that the qualitative
  presence and/or quantitation of that analyte in the sample cannot be determined with any degree of confidence.

Footnotes are applied to each qualifier to define the type of excursion that affected the sample result, resulting in the qualification of the data, as listed on this table.

Data are evaluated using the QA/QC criteria (including holding times and calibration) established in the applicable Quality Assurance Project Plan (QAPP), analytical methods and laboratory established control limits. Since the NJDEP validation guidelines apply to data generated using CLP methods, the application of these validation guidelines is modified to reflect method requirements, where applicable, since non-CLP methods are used in the analysis of samples.

A full QA/QC review is performed for 10 percent of the aqueous and solid samples, including a review of data summary forms and raw analytical data that were provided by the laboratory in the data package documentation. Partial review is performed for the remaining environmental samples submitted for data validation for this sampling event. Partial review consists of a review of the data summary forms. During the partial validation, only summary QA/QC forms are evaluated. The forms and the information contained on the forms are not evaluated for accuracy or completeness during the partial validation process.

The validation approach taken by O'Brien & Gere is a conservative one; qualifiers are applied to sample data to indicate both major and minor excursions. In this way, data associated with any type of excursion are identified to the data user. Major excursions will result in data being rejected, indicating that the data are considered unusable for either quantitative or qualitative purposes. Minor excursions will result in sample data being qualified as approximate that are otherwise usable for quantitative or qualitative purposes.

Excursions are subdivided into excursions that are within the laboratory's control and those that are out of the laboratory's control. Excursions involving laboratory control sample recovery, calibration response, method blank excursions, low or high spike recovery due to inaccurate spiking solutions or poor instrument response, holding times, interpretation errors, and quantitation errors are within the control of the laboratory. Excursions resulting from matrix spike recovery, serial dilution recovery, surrogate, and internal standard performance due to matrix interference from the matrix of the samples are examples of those excursions that are not within the laboratory's control if the laboratory has followed proper method control procedures, including performing appropriate cleanup techniques.

### **Parameter Type**

### **Approach in Applying Data Validation Qualifiers**

### Sample collection information-Cooler Temperature

Results for samples submitted for organic and inorganic analyses impacted by cooler temperatures of greater than  $10^{\circ}$ C are noted in the report.\* Qualifiers are not applied to data.

Sample collection information-Percent Solids

Results for samples submitted for organic and inorganic analyses that are impacted by percent solids of 50 percent are noted in the report.\* Qualifiers are not applied to data.

VOCs by USEPA Method 8260B Calibration Evaluation VOC target analytes are evaluated using the criteria of 15%RSD or correlation coefficient criteria of 0.990 for initial calibration curves. Calibration verifications are evaluated using a criterion of 20%D for all target compounds. Initial calibrations and calibration verifications were also evaluated using the criterion of a RF value of greater than or equal to a value of 0.01 for ketones and 0.05 for the remaining target analytes. If analyzed, the second-source standard (ICV) is evaluated using laboratory control limits or 70% to 130% recovery.

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Table 3 - O'Brien & Ger	e data validation approach using NJDEP data validation guidelines
VOCs by USEPA	VOC target analytes are evaluated using the criteria of 20%RSD or correlation coefficient criteria of 0.990 for initial calibration
Method 524.2	curves. Calibration verifications are evaluated using a criterion of 30%D for all target compounds. Initial calibrations and
Calibration Evaluation	calibration verifications were also evaluated using the criterion of a RF value of greater than or equal to a value of 0.05. If
	analyzed, the second-source standard (ICV) is evaluated using laboratory control limits or 70% to 130% recovery.
VOCs by USEPA	VOC target analytes are evaluated using the criteria of 35 percent relative standard deviation (%RSD) or
Method 624	correlation coefficient criteria of 0.990 for initial calibration curves. Calibration verifications are evaluated using
Calibration	criteria presented in Table 5 of USEPA Method 624 and 50 percent difference (%D) for the remaining target
Evaluation	analytes not listed in the method. Initial calibrations and calibration verifications are also evaluated using a
21010011	response factor (RF) criteria of greater than or equal to 0.05 for target analytes. A minimum of a RF pf 0.01 is
	required for ketones and poor-purging analytes. If analyzed, the second-source standard or low standard is
	evaluated using a 30% recovery or the laboratory control limits.
CVOCa by LICEDA	SVOC target analytes are evaluated using the criteria of 15 %RSD or correlation coefficient criteria of 0.990 for initial
SVOCs by USEPA Method 8270C	calibration curves. Calibration verifications are evaluated using a criterion of 20%D for all target compounds. Initial
Calibration Evaluation	calibrations and calibration verifications were also evaluated using the criterion of a RF value of greater than or equal to a
Calibration Evaluation	value of 0.05 for the target analytes. If analyzed, the second-source standard (ICV) is evaluated using laboratory control limits
	or 70% to 130% recovery.
	Due to any relative standard deviation (RSD) calibration excursions, detected results for analytes in samples associated with
	the calibration are qualified as approximate (J). Non-detected results associated with RSD excursions may be qualified as
	approximate (UJ) based on professional judgment.
Calibuation Astions for	If the RSD calibration excursion is greater than 90, detected results for analytes in samples associated with the calibration are
Calibration Actions for VOCs (8260B) and	qualified as approximate (J) and non-detected results may be rejected (R), applying professional judgment.
SVOCs (8270C)	Due to any %D calibration verification excursions, detected and non-detected results for analytes in samples associated with
37003 (82700)	the calibration are qualified as approximate (J, UJ).
	If the %D calibration excursion is greater than 90, detected results for analytes in samples associated with the calibration are
	qualified as approximate (J) and non-detected results may be rejected (R), applying professional judgment.
	For response factor excursions, detected results are qualified as approximate (J) and non-detected results are rejected (R).
	For initial calibration verifications (ICV) excursions, detected and non-detected results for analytes in samples associated with
	the calibration are qualified as approximate (J, UJ). The response direction and detection of target analytes in associated
PCBs by USEPA	sample may be considered in applying qualifiers.  PCB target analytes are evaluated using the criteria of 20 %RSD or correlation coefficient of 0.990 for initial calibration curves.
Method 8082	Calibration verifications are evaluated using a criterion of 15 %D for target analytes.
Calibration Evaluation	ICV recoveries are evaluated using a boratory control limits if available or 70 to 130%.
	Pesticide target analytes are evaluated using the criteria of 20 %RSD or correlation coefficient of 0.990 for initial calibration
Pesticides by USEPA	curves.
Method 8081A	Calibration verifications are evaluated using a criterion of 20 %D for the target analytes.
Calibration Evaluation	ICV recoveries are evaluated using laboratory control limits if available or 70 to 130%.
Harbisidas by HCEDA	Herbicide target analytes are evaluated using the criteria of 20 %RSD or correlation coefficient of 0.990 for initial calibration
Herbicides by USEPA Method 8151A	curves.
Calibration Evaluation	Calibration verifications are evaluated using a criterion of 20 %D for the target analytes.
Calibration Evaluation	ICV recoveries are evaluated using laboratory control limits if available or 70 to 130%.
	Due to any relative standard deviation (RSD) calibration excursions, detected results for analytes in samples associated with
	the calibration are qualified as approximate (J). Non-detected results associated with RSD excursions may be qualified as
Calibration Actions for	approximate (UJ) based on professional judgment.
PCB, Pesticides and	Due to any %D calibration verification excursions, detected and non-detected results for analytes in samples associated with
Herbicides GC analyses	the calibration are qualified as approximate (J, UJ).  For initial calibration verifications (ICV) excursions, detected and non-detected results for analytics in camples associated with
	For initial calibration verifications (ICV) excursions, detected and non-detected results for analytes in samples associated with
	the calibration are qualified as approximate (J, UJ). The response direction and detection of target analytes in associated sample may be considered in applying qualifiers.
Calibration Data- GC	Data are evaluated using the criteria of 20%RSD for initial calibrations, or correlation coefficient of 0.990 for calibration curves,
by USEPA Method	and 20%D for the calibration verifications. Results are qualified for primary column calibration excursions. The second-source
8011	standard (ICV) is evaluated using laboratory control limits or 70% to 130% recovery.
Organic Multi-results	When two results are reported, due to re-preparation or for dilution analyses, both sets of results are evaluated during the
J	validation process. Based on the evaluation of the associated quality control data, the results reflecting the higher quality data
	are reported.



Table 3 - O'Brien & Ger	e data validation approach using NJDEP data validation guidelines
General Organic Surrogate, MS/MSD,	Laboratory established control limits are used to assess duplicate, surrogate, MS/MSD, and LCS data.
Surrogate, MS/MSD, LCS, Duplicate Data	In the case that excursions are identified in more than one quality control sample of the same matrix within one sample delivery group, samples are batched according to sample preparation or analysis date and qualified accordingly.
	For surrogate recoveries are not within laboratory control limits:
	If two or more surrogate recoveries are outside of laboratory control limits for SVOC analysis, results are rejected (R, 81) unless matrix interferences are confirmed by re-extraction and reanalysis.
	If one or more surrogate recoveries are not within laboratory control limits for PCB, results are qualified as UJ, J, 81B).
	If LCS percent recoveries are less than laboratory control limits but greater than ten percent, non-detected and detected results are qualified as approximate (UJ, J, 88) to indicate minor excursions.
	If LCS percent recoveries are greater than laboratory control limits, detected results are qualified as approximate (J, 88) to indicate minor excursions.
	If LCS percent recoveries are outside of laboratory control limits and less than ten percent, detected results are qualified as approximate (J, 88) and non-detected results are qualified as rejected (R, 88A) to indicate major excursions.
	If RPDs for MSDs or duplicates are outside of laboratory control limits, detected results are qualified as approximate (J, 89A) to indicate minor excursions.
Organic MS/MSD Data	Qualification of organic data for MS/MSD analyses is performed only when both MS and MSD percent recoveries are outside of laboratory control limits with zero percent recovery.
	Organic data are rejected (R, 87) to indicate major excursions in the case that both MS/MSD recoveries are zero.
Sample dilution Data	Qualification of data is not performed if MS/MSD or surrogate recoveries are outside of laboratory control limits due to sample dilution.
MS/MSD and Field Duplicate Data – Organic Data	Qualification of data associated with MS/MSD or field duplicate excursions is limited to the un-spiked sample or the field duplicate pair, respectively.
Field Duplicate Data	Field duplicate data are evaluated against relative percent difference (RPD) criteria of less than 50 percent for aqueous samples and less than 100 percent for soils when results are greater than five times the QL. When sample results for field duplicate pairs are less than five times the QL, the data are evaluated using control limits of plus or minus two times the QL for soils. If RPDs for field duplicates are outside of laboratory control limits, detected and non-detected results are qualified as approximate (UJ, J, 90) to indicate minor excursions.
Internal Standard - Organic Data	Internal standard recoveries are evaluated using control limits of within 50% of the lower standard area and up to 100% of the upper standard area of the associated calibration verification standard.
	Sample results are qualified as approximate (UJ, J, 50) if one internal standard does not meet criteria.
	Detected sample results are qualified as approximate (J, 51) if two or more internal standards do not meet criteria.
Internal	Non-detected sample results are rejected (R, 51) if two or more internal standards do not meet criteria.
Standard/Surrogate - Organic Data- Drinking Water methods	Internal standard recoveries are evaluated using method control limits. Monitor the integrated areas of the quantitation ions of the internal standards and surrogates in all samples, continuing calibration checks, and blanks. These should remain reasonably constant over time. An abrupt change may indicate a matrix effect or an instrument problem. If a cryogenic interface is utilized, it may indicate an inefficient transfer from the trap to the column. These samples must be reanalyzed or a laboratory fortified duplicate sample analyzed to test for matrix effect. A drift of more than 50% in any area is indicative of a loss in sensitivity, and the problem must be found and corrected.
	CCV- Determine that the absolute areas of the quantitation ions of the internal standard and surrogates have not decreased by more than 30% from the areas measured in the most recent continuing calibration check, or by more than 50% from the areas measured during initial calibration. If these areas have decreased by more than these amounts, adjustments must be made to restore system sensitivity.
Evaluation of Internal	Internal standard areas of samples are evaluated using the validation control limit of 70 to 130 percent recovery when compared to the calibration verification associated with the samples.
Standards for samples (VOCs for USEPA	Sample results are qualified as approximate (UJ, J, 50) if one internal standard does not meet criteria.
Method 524.2)	Detected sample results are qualified as approximate (J, 51) if two or more internal standards do not meet criteria.
	Non-detected sample results are rejected (R, 51) if two or more internal standards do not meet criteria.



Table 3 - O'Brien & G	Gere data validation approach using NJDEP data validation guidelines
	Internal standard areas of CCVs are evaluated using the validation control limit of 50 to 100 percent recovery when compared to the initial calibration.
<b>Evaluation of CCVs</b>	
(VOCs for USEPA	Sample results are qualified as approximate (UJ, J, 50) if one internal standard does not meet criteria.
Method 524.2)	Detected sample results are qualified as approximate (J, 51) if two or more internal standards do not meet criteria.
	Non-detected sample results are rejected (R, 51) if two or more internal standards do not meet criteria.
Evaluation of Initial	
(ICV) and Calibration	
Verification (CCV) fo	
Metals by 6010B/6020A,	Mercury is evaluated using the criteria for ICV of 90% to 110% of the expected value and 80% to 120% of the expected value for the CCV.
Mercury by	Total Cyanide is evaluated using the criteria for ICV and CCV of 85% to 115% of the expected value.
7470A/7471B, and	For analyses utilizing a calibration curve, the correlation coefficient for the first or second order curve must be $\geq 0.995$ .
Total Cyanide by	
9012B	
Performance	ICP-MS data is evaluated using resolution of mass calibration of within 0.1 $\mu$ and the %RSD of less than 15%.
Evaluation for ICP-M	Resolution must be less than 0.9amu of full width at 10% of peak height.
by 6020A	
Evaluation of Initial (ICV) and Calibration	
Verification (CCV) for	
Metals by EPA metho	
200.7/200.8 and	For analyses utilizing a calibration curve, the correlation coefficient for the first or second order curve must be $\geq 0.995$ .
Anions by Method	
300.0	
Evaluation of Internal	
Standards for ICP-MS	response in the calibration blank.
by 200.8	Internal standard recognition are such at all views and tall limits of a consent relative internal to (0/DI) from COO/ to 1250/ of the
Evaluation of Internal Standards for ICP-MS	
by 6020A	The intensity of any internal standard must be >30% or <120% of the intensity of the internal standard in the initial calibration
by 6626/1	standard.
	The intensity of the internal standard of the CCB and CCV must agree within ±20% of the intensity of the internal standard in
	the ICV.
Metal and Inorganic	
MS/MSD,	samples for the same matrix, within the same preparation batch, within the same SDG group.
Laboratory/Field	
Duplicate, Serial Dilution	
Validation Footnotes	
Footnote	Type of Excursion
1	The value reported is less than or equal to three (3) times the value in the method blank/preparation blank. It is the policy of NJDEP-
	DPFSR to negate the reported value due to probable foreign contamination unrelated to the actual sample. The end-user, however, is
	alerted that a reportable quantity of the analyte/compound was detected. The B qualifier must be reported.
2	The value reported is greater than three (3) times but less than or equal to 10 times the value in the method blank/preparation blank
	and is considered "real". However, the reported value must be quantitatively qualified "J" due to the method blank contamination. The
3	"B" qualifier alerts the end-user to the presence of this analyte/compound in the method blank.  The value reported is less than or equal to three (3) times the value in the trip/field blank. It is the policy of NJDEP-DPFSR to negate the
3	reported value as due to probable foreign contamination unrelated to the actual sample. The end-user, however, is alerted that a
	reportable quantity of the analyte/compound was detected.
4	The value reported is greater than three (3) times the value in the trip/field blank but less than or equal to 10 times the value in the
	blank and is considered "real". However, the reported value must be quantitatively qualified "J" due to trip/field blank contamination.
4A	The result was qualified due to negative drift.
4B	The result was qualified as "U" due to blank contamination.
5	The concentration reported by the laboratory is incorrectly calculated.
6	The laboratory failed to report the presence of the analyte in the sample.
7	The reported metal value was qualified because the Initial/Continuing Calibration Standard was not within the recovery range.
8	No CRDL Standard for AA or ICP analysis was performed. Therefore, the analyte affected was rejected.
9	The reported concentration was quantitatively qualified because the concentration was below the CRDL but greater than the MDL. The
	concentration is considered estimated since the value obtained is at the low end of the instrument performance.
9A	IDLs are greater than the CRDLs.



Table 3 - O'Bri	en & Gere data validation approach using NJDEP data validation guidelines			
10	The reported metal value was qualified because the ICP Interference Check Sample was outside the recovery range (80-120 percent).			
11	The non-detect metal value was qualified "UJ" because the ICP Interference Check Sample was within the range of 50 and 79%; hence a possibility of false negatives exists.			
12	This non-detected metal analyte had Laboratory Control Sample recovery that fell within the range of 70-79%. The end-user should be aware of the possibility of false negatives; therefore, this analyte is flagged as estimated (UJ).			
13	The reported metal value was qualified because the Laboratory Control Sample recovery fell within the range of 70-79 %. The er user should be aware of results that may be biased low.			
14	The reported metal value was qualified because the Laboratory Control Sample recovery was greater than 120% but less than or equa to 130%. The end-user should be aware of results that may be biased high.			
15	The metal analyte is rejected because the Laboratory Control Sample recovery was less than 70% or greater than 130%.			
16	In the Duplicate Sample Analysis for metals, the analyte fell outside the control limits of +20 percent or + CRDL. Therefore, result for the metal was qualified.			
17	This analyte was rejected because the laboratory performed the Duplicate Analysis on a field blank.			
18	The reported metal value was qualified because the spike recovery was greater than 125 percent but less than or equal to 200%.			
18A	The reported metal was qualified because both the spike recovery and matrix spike duplicate recovery were outside of the validation control limits.			
19	The reported metal value was qualified because the spike recovery was between 25 and 74 percent.			
20	The reported metal value was qualified because the spike recovery was less than 25 percent. The reported value actually indicated the minimum concentration at which the metal was present.			
21	The non-detected metal value was qualified (UJ) because the spike recovery was between 25 and 74 percent. The possibility of a false negative exists.			
22	The non-detected metal value was rejected because the spike recovery was less than 25 percent.			
23	The reported metal value was rejected because the laboratory used a field blank for the Sample Spike Analysis.			
24	There was no Post-Digestion Spike Sample Recovery analysis performed. Therefore, the analyte was rejected.			
25	The reported metal value was qualified because the Serial Dilution was not within ten percent of sample concentration.			
26	The reported metal value was rejected because the laboratory used a field blank for the Serial Dilution analysis or the post-diges spike.			
27	This metal analyte is rejected because the preparation blank concentration of this analyte is greater than the CRDL and the reported sample concentration is less than ten (10) times the preparation blank concentration.			
28	The laboratory incorrectly transcribed the raw data onto the Inorganic Analysis Data Sheet form or there are data package issues.			
28A	Verification of instrument parameters was performed outside of the required frequency.			
28B	A percent solids issue was detected.			
29	The reported metal analyte was rejected because the CRDL standard % Recovery fell less than 30% or was greater than 175%, or another severe CRDL deficiency was detected.			
30	The non-detected metal value was rejected because the post-digestion spike recovery was less than 25 percent.			
30A	The metal value was qualified since the post-digestion spike recovery was exceeded.			
31	The reported metal analyte was rejected because the associated Continuing Calibration Blank result was greater than the CRDL.			
32	The reported metal analyte was rejected because this sample is not associated with a Laboratory Control Sample or ICB or CCB.			
33	The laboratory made a transcription error.			
33A	A methods comparison issue was detected.			
34	The laboratory used an incorrectly associated Preparation Blank.			
35	This analyte is rejected because the laboratory exceeded the holding time for analysis or extraction.			
35A	Result was qualified due to a holding time excursion.			
36	This metal value was qualified because the CRDL standard was not within the recovery range.			
37	The reported concentration is quantitatively qualified due to calibration deficiencies.			
38	The reported concentration is quantitatively qualified due to surrogate recovery outliers.			
39	The reported concentration is quantitative qualified because the concentration is below the RL.			
40	The sample holding time to re-extraction and/or reanalysis was exceeded. All positive results including the tentatively identified compounds are highly qualified.			
41	The mass spectral identification has not been confirmed and the identification of this compound has been rejected. This compound should now be considered an unknown and the reported concentration is considered an estimated value.			
42	The percent Difference of the calculated values on both columns is greater than 100% and less than 999.9 %. This value is significantly greater than the 25 % limits established by the USEPA-Contract Laboratory Program. The extreme variation between the values from the two columns is apparently due to instrumentation problems and/or matrix interference. Therefore, the reported concentrations cannot be verified and only a tentative identification of the Aroclor or pesticide can be determined.			
42A	The percent difference from both columns was greater than 25%.			
42B	The percent difference from both columns was greater than 40%.			



Table 3 - O'Brien	& Gere data validation approach using NJDEP data validation guidelines	
42C	The percent difference from both columns was greater than 70%.	
42D	The percent difference from both columns was greater than 100% without evidence of matrix interferences being present. The results are rejected (R).	
42E	Results were reported at a concentration that was less than the PQL with a %D greater than 50 percent. The PQL is reported and qualified as non-detected (U).	
43	The peak retention times of the Aroclors or pesticides detected in the samples are outside of the retention time window established in the initial calibration. The identification of the Aroclors or pesticides cannot be verified due to the retention time shift outside of the windows. Retention time shifts are evident in all of the continuing calibration standards and the Performance Evaluation Mixtures,	
	therefore the usability of the data is questionable.	
44	The laboratory didn't provide the mass spectral proof for the analyte although the quantitation report indicates the presence of the analyte. The presence of this analyte in the sample is considered tentative.	
45	The non target compound is qualified "J" and considered an estimated value because relative response factors are not determin non-target compounds.	
46	The laboratory's call on the non target compound did not match the mass spectra of the compound at the approximate scan number in the blank. The laboratory call is incorrect.	
47	The laboratory failed to report this analyte on the Organic Analysis Data Sheet (OADS)  Form even though the TIC, quantitation report and library search indicates a hit for the analyte.	
48	The laboratory reported this analyte in the QADS form. However, this analyte was negated in the quantitation report. QA reviewer agrees the mass spectrum is not a good match and therefore, negates the presence of this analyte in the sample.	
49	No library search was submitted for this unknown.	
49A	Results were rejected since correct internal standard was not used.	
50	One internal standard area in the sample did not meet the QC criteria. Therefore, all compound results using this internal standard for quantitation are quantitatively estimated. (UJ, J)	
51 (See 84)	Two or more internal standard areas in the sample did not meet the QC criteria with recoveries of greater than 25%. The detected results for the entire fraction for that sample are qualified as approximate (J). The non-detected results are rejected (R).	
52	The RIC in the raw data indicates a non-target(s) is present. The lab failed to report and provide library search(s) for the non-target(s).	
53	The laboratory did not quantify the pesticides present in the sample. The pesticide was confirmed on a second column. Quantitation of the peaks revealed that the value is above the CRQL.	
54	The lab failed to report this analyte although it was found in both columns and is within the retention times of both columns for the analyte.	
55	The retention time window for this analyte overlaps with the retention time window of another analyte. The identity is indistinguishable and therefore tentative.	
56	The laboratory reported concentration does not agree with QA reviewer's calculated concentration.	
57	The compound exceeded the calibration range of the instrument and is indicated with the "E" qualifier.	
58	The compound is a suspected Aldol condensation product and is flagged with the "A" qualifier.	
59	The laboratory was required to dilute the samples to bring the peaks onto scale.	
60	This sample was diluted prior to analysis. The value reported prior to the dilution correction is less than three (3) times the value in the method blank. It is the policy of NJDEP-DPFSR to negate the reported value due to probable foreign laboratory contamination unrelated to the actual sample. The end-user is alerted that a reportable quantity of the analyte was detected.	
61	This non-target compound was detected as a target compound in another analytical fraction. Therefore, the presence of this compound as a non-target analyte is negated.	
62	This sample was diluted prior to analysis. The value reported prior to the dilution correction is greater than three (3) times the value in the method blank and is considered "real". However, the reported value must be quantitatively qualified "J" due to method blank contamination. The "B" qualifier alerts the end-user to the presence of this analyte in the method blank.	
62A	Results are rejected due to a severe blank analysis excursion.	
62B	Results are qualified due to a blank analysis excursion.	
63	The results are rejected because the initial calibration, continuing calibration or internal standard was not performed using the proper sequence, concentration, matrix, or internal standards.	
63A	Results are rejected due to a severe pesticide/Aroclor analysis issue.	
63B	Results are negated due to a blank analysis excursion.	
63C	Results are qualified due to a pesticide/Aroclor analysis issue.	
64	The results are rejected because the D of the single component pesticide and/or surrogate in the PEM(s) is greater than 25%.	
64A	Results are rejected due to a major calibration excursion.	
65	The results are rejected because of resolution, scaling, or retention time issues.	
65A	Results are qualified due to scaling, or calibration issues.	
66	The result is rejected due to retention time deficiencies.	
67	The result is qualified because the DDT and/or Endrin breakdown was greater than 20%.	



Table 3 - O'Bri	en & Gere data validation approach using NJDEP data validation guidelines	
68	The result is qualified because the combined DDT/Endrin breakdown is greater than	
	30%.	
69	The results are rejected because GPC cleanup was not performed on the sample extract.	
70	The results are rejected because florisil cleanup was not performed on the sample extract.	
71	The results are rejected due to GPC calibration or analysis deficiencies.	
72	The results are rejected because the florisil cartridge check yielded unacceptable percent recoveries or was not performed	
	properly.	
73	The sample holding time was exceeded by greater than ten days. The sample results are rejected.	
74	The GC/MS Instrument Performance Check Solution (IPCS) failed acceptance criteria or was not performed. The associ sample results are rejected.	
74A	The results are qualified due to IPCS time-of-analysis excursions.	
75	Three or more analytes in the initial calibration or continuing calibration failed to meet acceptance criteria. The associated sample results are rejected.	
76	The results in the fraction are rejected because the response factor in the initial and/or continuing calibration is less than 0.01 or does not meet the project requirement.	
77	The results in the fraction are rejected because the %RSD and/or %D is greater than 40% (or in the case of %D, less than - 40%).	
78	The positive result is qualified because the RRF of the compound (with no %RSD or %D) is less than 0.01 or does not meet the project requirement.	
79	The non-detect result is rejected because the RRF of the compound (with no %RSD or %D) is less than 0.01.	
80	Five or more analytes in the initial calibration or continuing calibration failed to meet %RSD or %D and/or RRF acceptance criteria. The associated sample results are rejected.	
80A	Results are rejected since the continuing calibration was not performed properly.	
81	Sample results for the fraction are rejected because the % recovery of two or more SMCs (or surrogates) failed to meet criteria.	
81A	Results are rejected due to severe surrogate analysis excursions.	
81B	Results are qualified due to surrogate analysis excursion.	
82	Sample results for the fraction are rejected because the %recovery of one or more SMCs (or surrogates) in the associated method blank failed to meet criteria.	
83	Sample results for the fraction are rejected because the retention time of one or more internal standards deviated by more than +/-30 seconds from the retention time of the corresponding internal standard in the associated calibration standard.	
84	Two or more internal standard areas in the sample did not meet the QC criteria with recoveries of less than 25%. The detected results and non-detected results are rejected (R).	
84A	Results are qualified due to sulfur cleanup issue.	
84B	Results are qualified due to internal standard failure.	
85	Sample results for the fraction are rejected because sulfur was present in the sample and sulfur cleanup was not performed or performed properly.	
86	Results are rejected due to failure to submit manual integration technique.	
87	Results are rejected or qualified due to zero matrix spike/ matrix spike duplicate recoveries.	
88	Results are qualified due to laboratory control sample excursions.	
88A	Results are rejected due to laboratory control sample recoveries of less than ten percent.	
89	Detected organic results are qualified due to zero matrix spike/matrix spike duplicate recoveries.	
89A	Organic results are qualified due to matrix spike/matrix spike duplicate precision excursions.	
90	Results are qualified due to field duplicate excursions. (UJ, J)	
91	Results are qualified due to calibration excursions.	
92	Results are rejected due to significant canister pressure differences.	
93	Results are rejected since SIM was utilized.	
94	Results are rejected since a separate MDL study was not performed for each instrument.	
95	Results are qualified due to analysis excursions.	
96	Results are qualified due to a sample collection excursion.	
96A	Results are rejected due to a sample collection excursion.	
97	Results are qualified due to sample preparation excursion.	
98	The reported hexavalent chromium result was qualified because the post verification spike was greater than 115%.	
99	The reported hexavalent chromium result was qualified because the post verification spike was less than 85%	
100	The non-detected hexavalent chromium result was qualified (UJ) because the post verification spike was less than 85%. The possibility	



Table 3 - O'Brien & Gere data validation approach using NJDEP data validation guidelines		
	of a false negative exists.	
101	The reported hexavalent chromium result was qualified because the pre-digestion spike recovery was less than 75%.	
102	The reported hexavalent chromium result was qualified because the pre-digestion spike recovery was greater than 125%.	
103	The non-detected hexavalent chromium result was qualified because the pre-digestion spike recovery was less than 75%. The possibility of a false negative exists.	
104	Results are qualified due to sample preservation excursion.	
* Indicates t	hat NJDEP data validation guidelines do not address this situation; therefore, validation qualifiers are not applied to data.	
Source O'Bri	en & Gere	

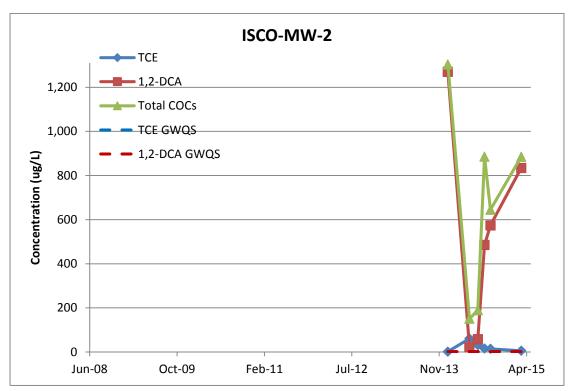
QA/QC Term	Definition
QA/QC Term	The level above which numerical results may be obtained with a specified degree of
Quantitation limit	confidence; the minimum concentration of an analyte in a specific matrix that can be identified and quantified above the method detection limit and within specified limits of precision and bias during routine analytical operating conditions.
Method detection limit	The minimum concentration of an analyte that undergoes preparation similar to the environmental samples and can be reported with a stated level of confidence that the analyte concentration is greater than zero.
Instrument detection limit	The lowest concentration of a metal target analyte that, when directly inputted and processed on a specific analytical instrument, produces a signal/response that is statistically distinct from the signal/response arising from equipment "noise" alone.
Gas chromatography/mass spectrometry (GC/MS) instrument performance check	Performed to verify mass resolution, identification, and to some degree, instrument sensitivity. These criteria are not sample specific; conformance is determined using standard materials.
Calibration	Compliance requirements for satisfactory instrument calibration are established to verify that the instrument is capable of producing acceptable quantitative data. Initial calibration demonstrates that the instrument is capable of acceptable performance at the beginning of analysis and calibration verifications document satisfactory maintenance and adjustment of the instrument on a day-to-day basis.
Relative Response Factor	A measure of the relative mass spectral response of an analyte compared to its internal standard. Relative Response Factors are determined by analysis of standards and are used in the calculation of concentrations of analytes in samples.
Relative standard deviation	The standard deviation divided by the mean; a unit-free measure of variability.
Correlation coefficient	A measure of the strength of the relationship between two variables.
Relative Percent Difference	Used to compare two values; the relative percent difference is based on the mean of the two values, and is reported as an absolute value, i.e., always expressed as a positive number or zero.
Percent Difference	Used to compare two values; the percent difference indicates both the direction and the magnitude of the comparison, i.e., the percent difference may be either negative, positive, or zero.
Percent Recovery	The act of determining whether or not the methodology measures all of the target analytes contained in a sample.
Calibration blank	Consists of acids and reagent water used to prepare metal samples for analysis. This type of blank is analyzed to evaluate whether contamination is occurring during the preparation and analysis of the sample.
Method blank	A water or soil blank that undergoes the preparation procedures applied to a sample (i.e., extraction, digestion, clean-up). These samples are analyzed to examine whether sample preparation, clean-up, and analysis techniques result in sample contamination.
Field/equipment	Collected and submitted for laboratory analysis, where appropriate. Field/equipment blanks are handled in the same manner as environmental samples. Equipment/field blanks are analyzed to assess contamination introduced during field sampling procedures.
Trip blank	Consist of samples of analyte-free water that have undergone shipment from the sampling site to the laboratory in coolers with the environmental samples submitted for volatile organic compound (VOC) analysis. Trip blanks will be analyzed for VOCs to determine if contamination has taken place during sample handling and/or shipment. Trip blanks will be utilized at a frequency of one each per cooler sent to the laboratory for VOC analysis.
Internal standards performance	Compounds not found in environmental samples which are spiked into samples and quality control samples at the time of sample preparation for organic analyses. Internal standards must meet retention time and recovery criteria specified in the analytical method. Internal standards are used as the basis for quantitation of the target analytes.
Surrogate recovery	Compounds similar in nature to the target analytes but not expected to be detected in the environmental media which are spiked into environmental samples, blanks, and quality control samples prior to sample preparation for organic analyses. Surrogates are used to evaluate analytical efficiency by measuring recovery.

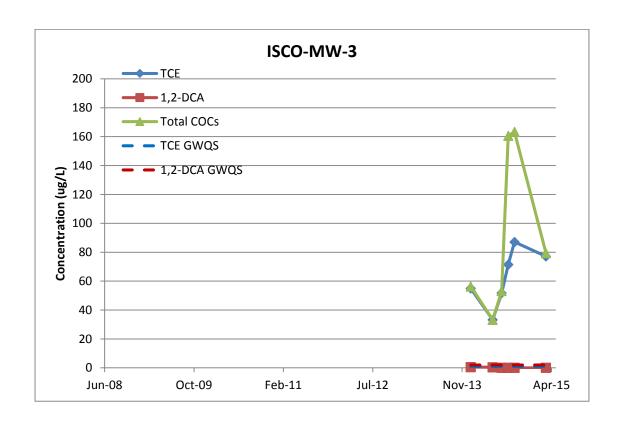


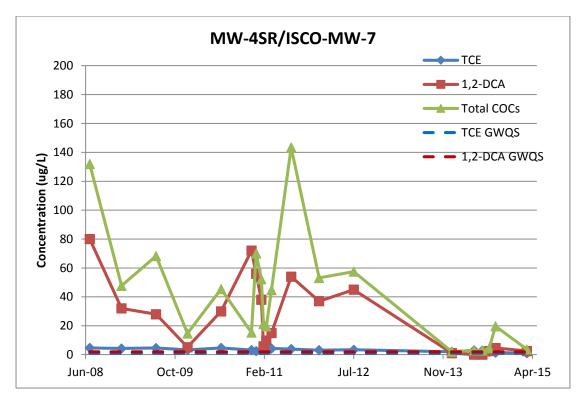
Laboratory control sample Matrix spike blank analyses	Standard solutions that consist of known concentrations of the target analytes spiked into laboratory analyte-free water or sand. They are prepared or purchased from a certified manufacturer from a source independent from the calibration standards to provide an independent verification of the calibration procedure. They are prepared and analyzed following the same procedures employed for environmental sample analysis to assess method accuracy independently of sample matrix effects.
Laboratory duplicate	Two or more representative portions taken from one homogeneous sample by the analyst and analyzed in the same laboratory.
Matrix	The material of which the sample is composed or the substrate containing the analyte of interest, such as drinking water, waste water, air, soil/sediment, biological material.
Matrix Spike (MS)	An aliquot of a matrix (water or soil) fortified (spiked) with known quantities of specific target analytes and subjected to the entire analytical procedure in order to indicate the appropriateness of the method for the matrix by measuring recovery.
Matrix spike duplicate (MSD)	A second aliquot of the same matrix as the matrix spike that is spiked in order to determine the precision of the method.
Retention time	The time a target analyte is retained on a GC column before elution. The identification of a target analyte is dependent on a target compound's retention time falling within the specified retention time window established for that compound.
Relative retention time	The ratio of the retention time of a compound to that of a standard.

# Attachment 4 Concentration Trend Graphs

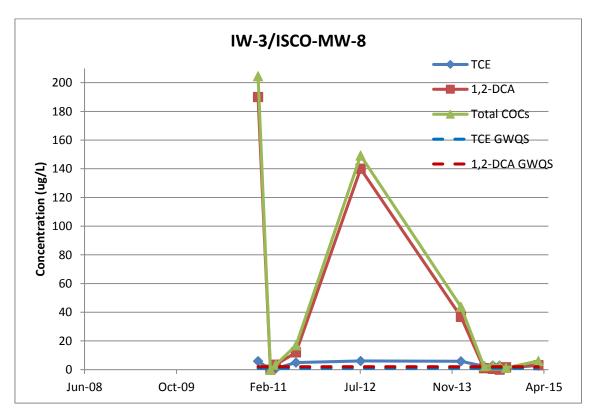
## **ISCO Treatment Area 1 Wells**





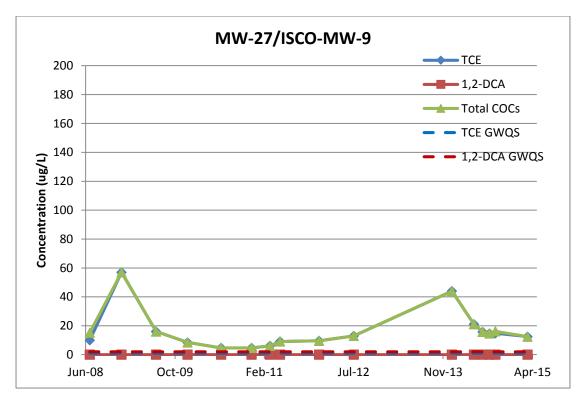


Note: ISCO-MW-7 was installed to replace former well MW-4SR on December 23, 2013, in accordance with the approved RDR/RAWP.

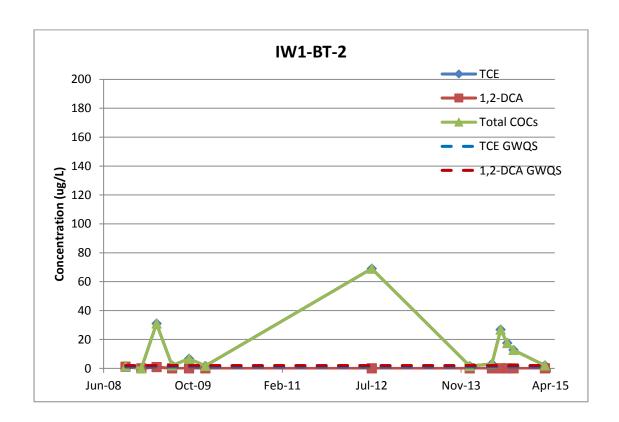


Note: ISCO-MW-8 was installed to replace former well IW-3 on November 27, 2013, in accordance with the approved RDR/RAWP.

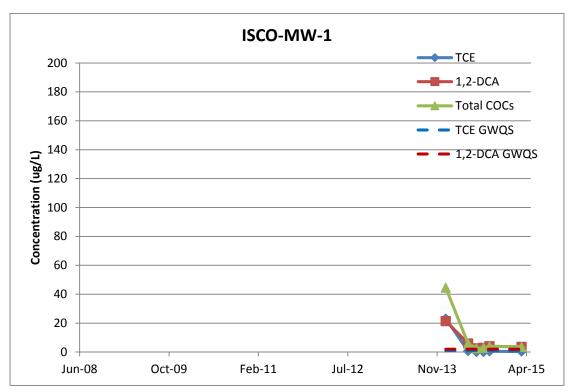


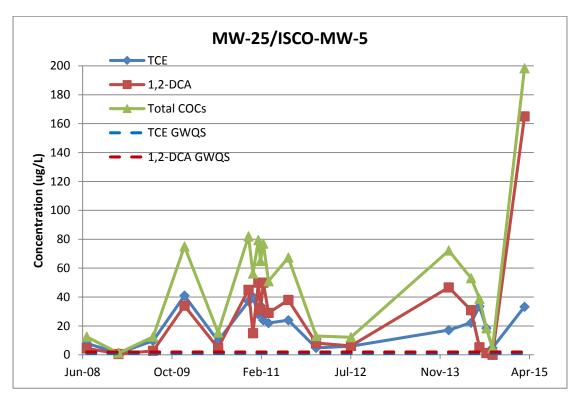


Note: ISCO-MW-9 was installed to replace former well MW-27 on December 23, 2013, in accordance with the approved RDR/RAWP.



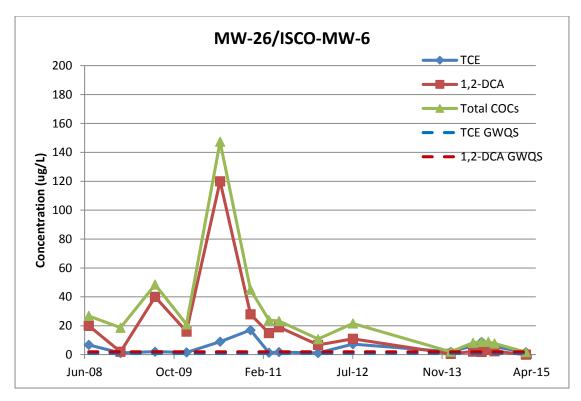
### **ISCO Treatment Area 2 Wells**



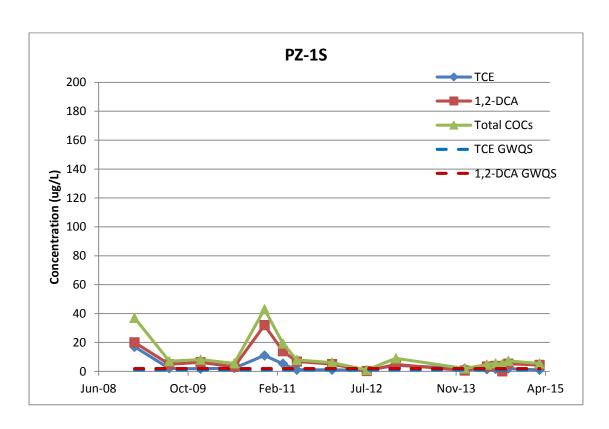


 $Note: ISCO-MW-5 \ was \ installed \ to \ replace \ former \ well \ MW-25 \ on \ November \ 26, 2013, in \ accordance \ with \ the \ approved \ RDR/RAWP.$ 

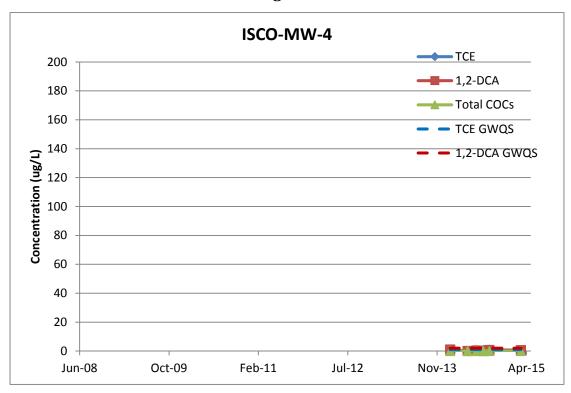


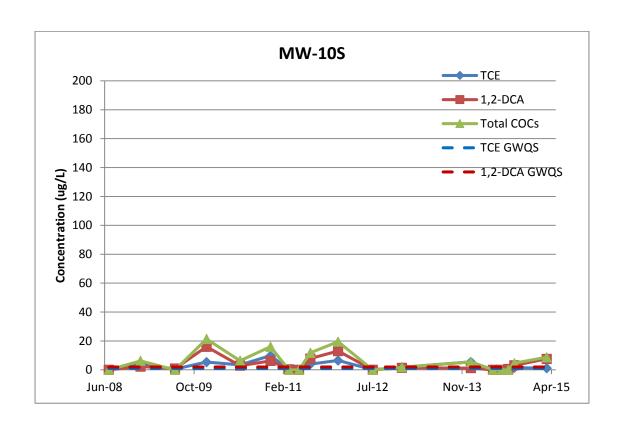


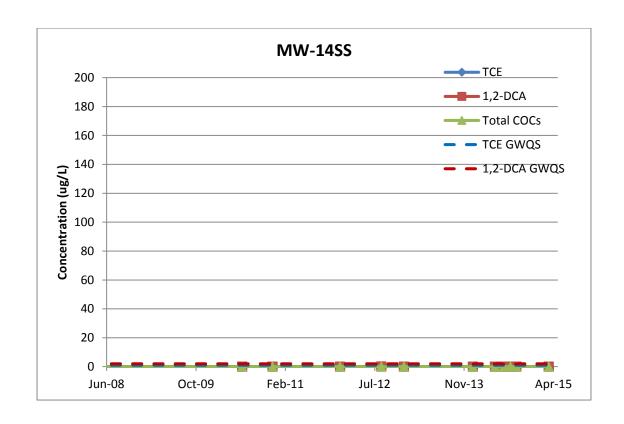
Note: ISCO-MW-6 was installed to replace former well MW-26 on November 27, 2013, in accordance with the approved RDR/RAWP.

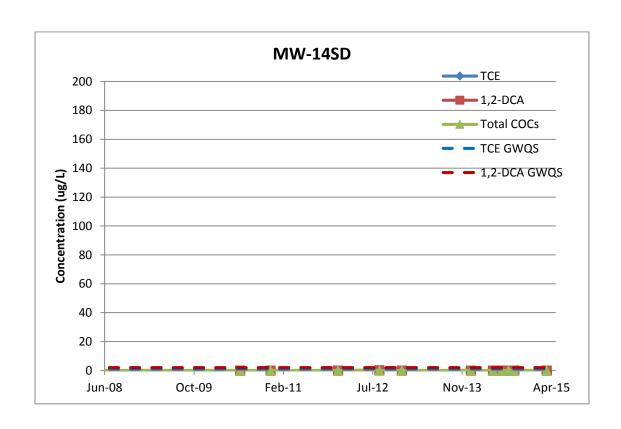


# **ISCO Downgradient Wells**

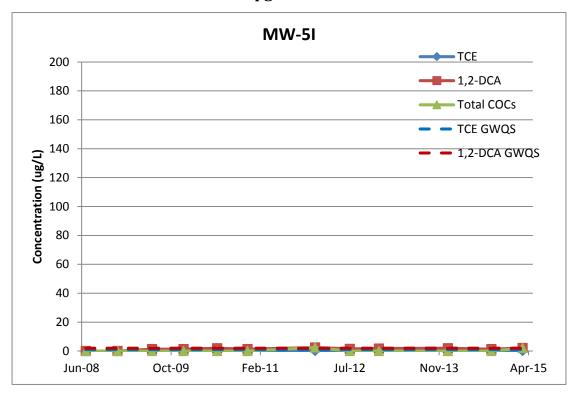


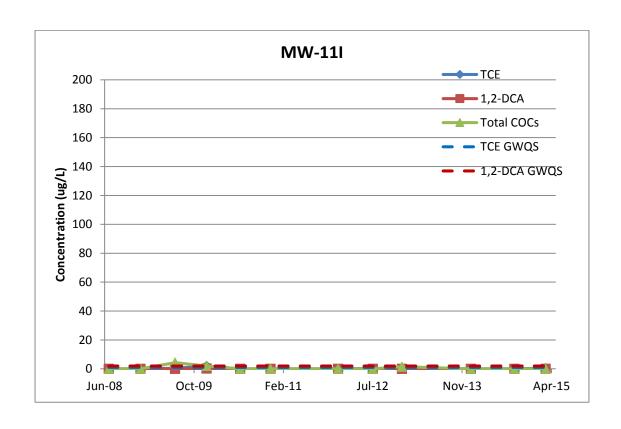


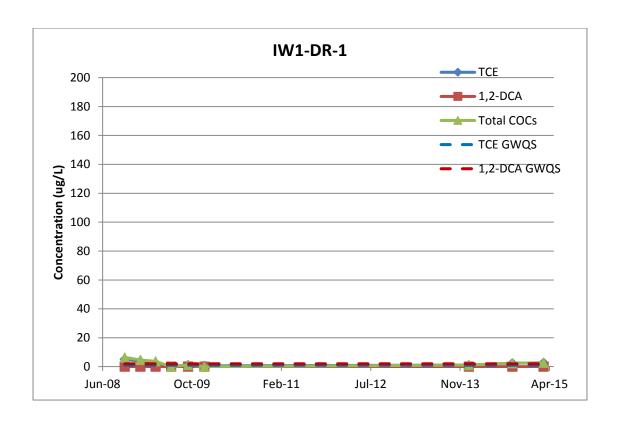




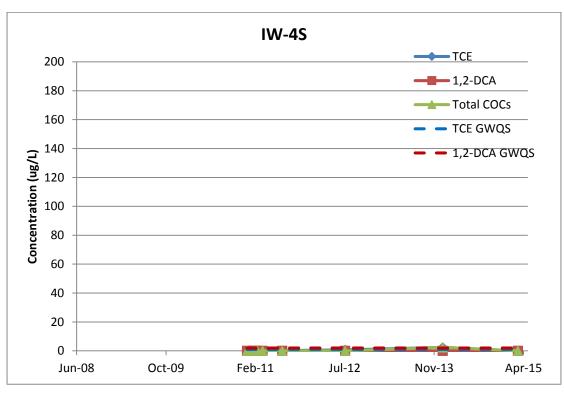
# **ISCO Upgradient Wells**

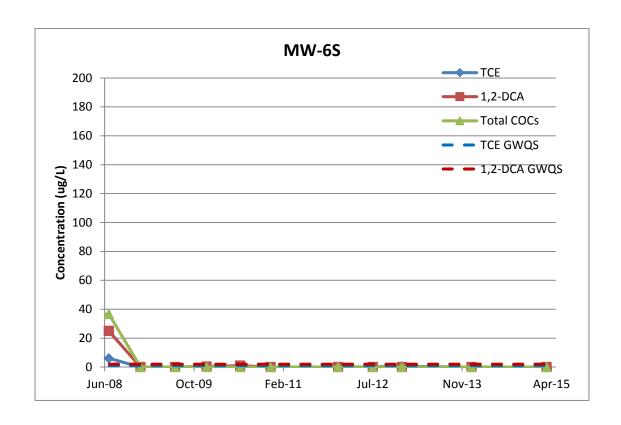


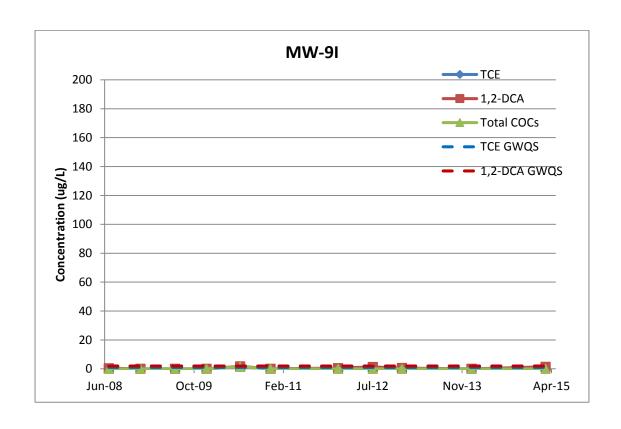


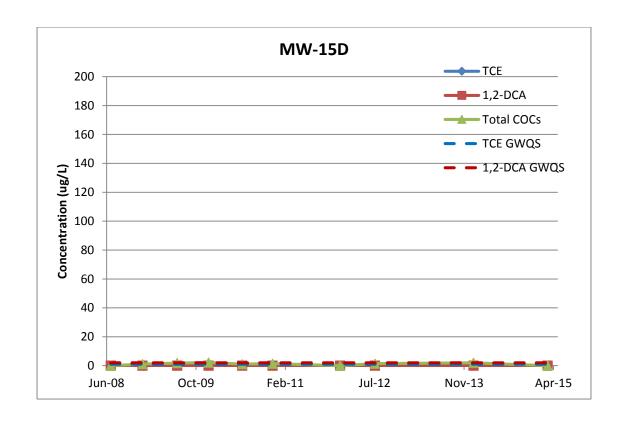


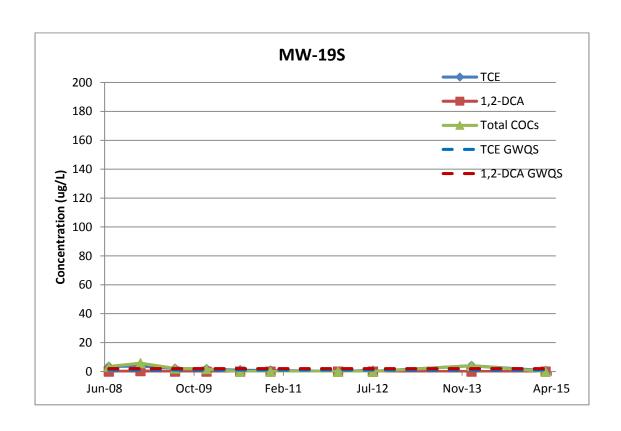
### Other

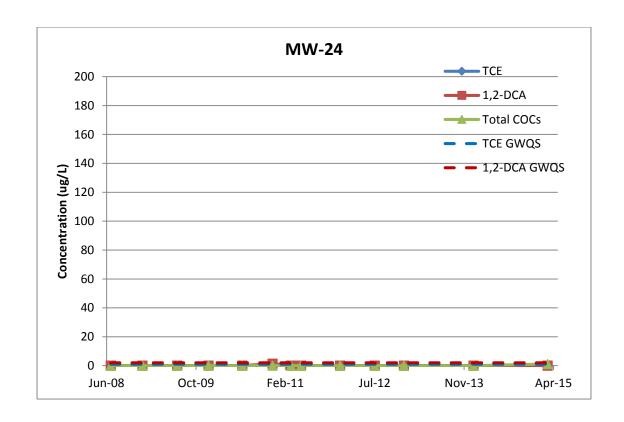


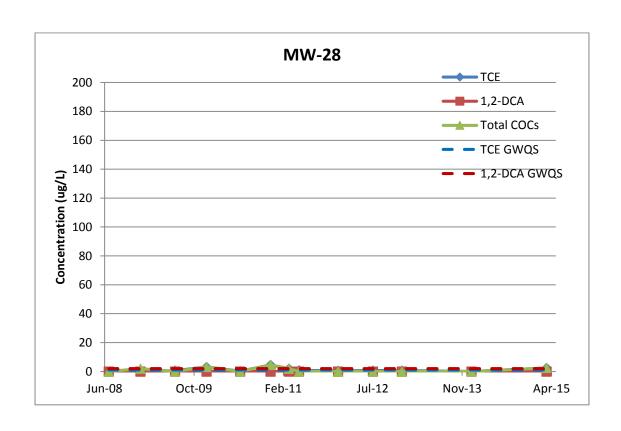


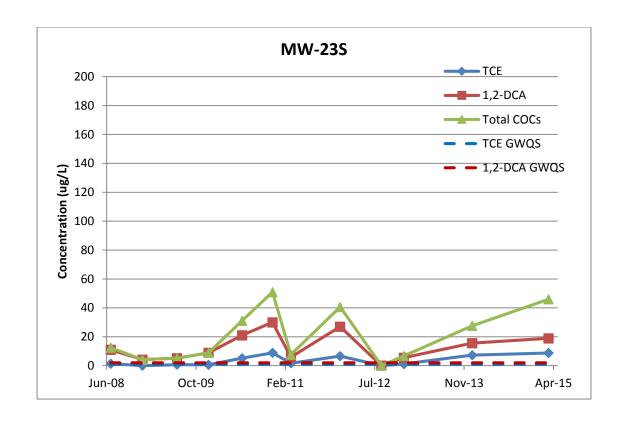


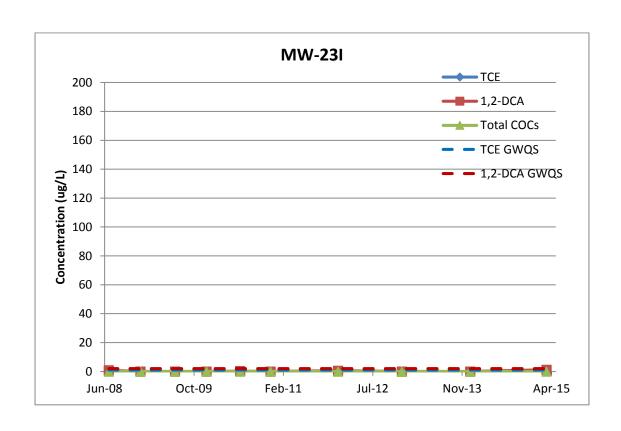


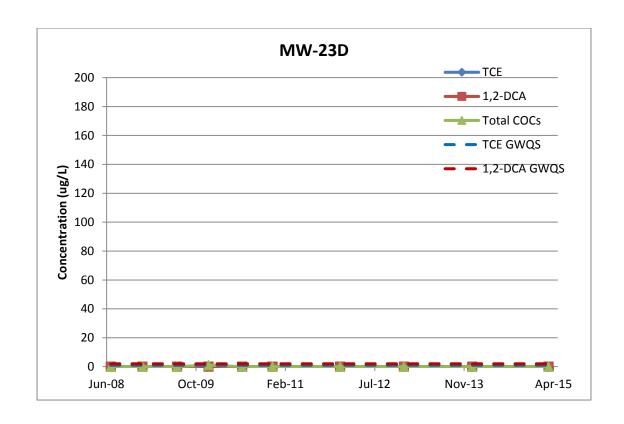


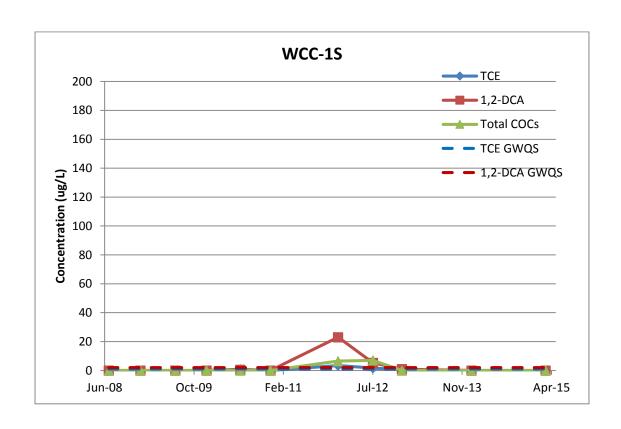


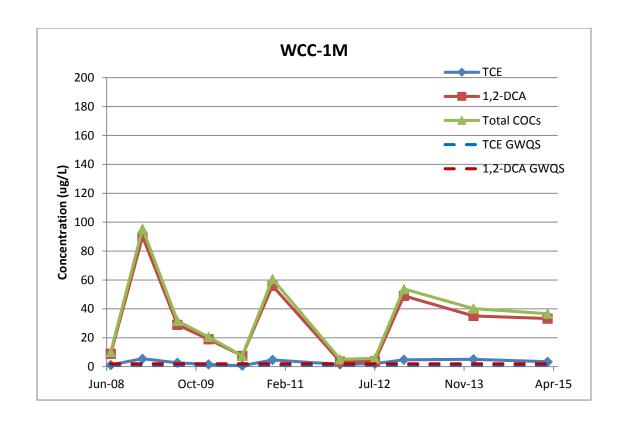


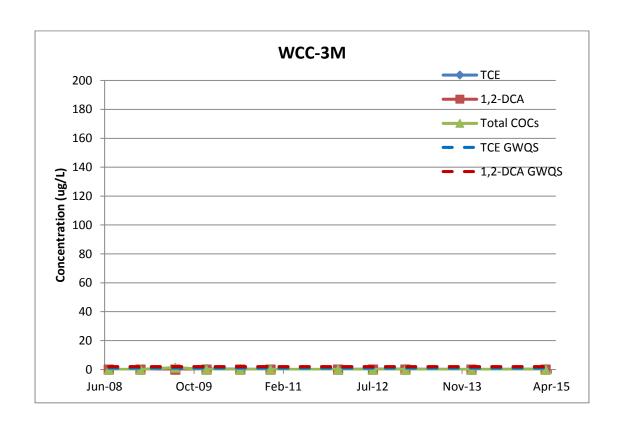


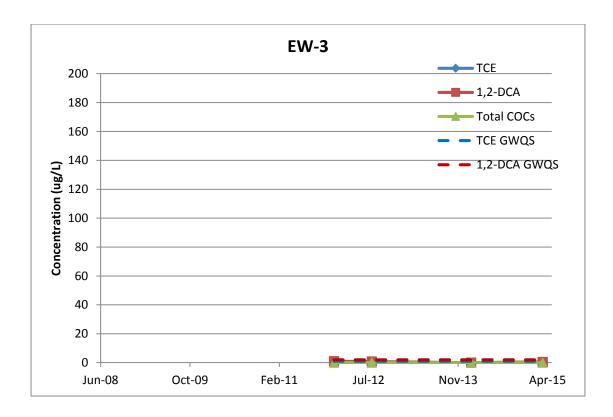












\*Total COCs represents the sum of those groundwater constituents above their respective New Jersey Groundwater Quality Standard (NJGWQS)